

Data Path : Z:\svoasrv\HPCHEM1\BNA\_E\Data\BE110824\  
 Data File : BE101544.D  
 Acq On : 8 Nov 2024 10:07  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

**Instrument :**  
 BNA\_E  
**ClientSampleId :**  
 SSTDCCC040

**Manual Integrations**  
**APPROVED**  
 Reviewed By :Yogesh Patel 11/10/2024  
 Supervised By :mohammad ahmed 11/11/2024

Quant Time: Nov 08 10:14:10 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_E\Methods\8270-BE110624.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 07 00:01:20 2024  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
<b>Internal Standards</b>							
1) 1,4-Dichlorobenzene-d4	7.558	152	51174	20.000	ng	0.00	
21) Naphthalene-d8	10.326	136	237509	20.000	ng	0.00	
39) Acenaphthene-d10	14.174	164	179651	20.000	ng	0.00	
64) Phenanthrene-d10	16.912	188	420948	20.000	ng	0.00	
76) Chrysene-d12	21.072	240	472751	20.000	ng	0.00	
86) Perylene-d12	23.357	264	602462	20.000	ng	0.00	
<b>System Monitoring Compounds</b>							
5) 2-Fluorophenol	5.314	112	236611	81.754	ng	0.00	
7) Phenol-d6	6.941	99	328691	82.833	ng	0.00	
23) Nitrobenzene-d5	8.769	82	312679	83.156	ng	0.00	
42) 2,4,6-Tribromophenol	15.678	330	273256	80.833	ng	0.00	
45) 2-Fluorobiphenyl	12.793	172	856759	77.869	ng	0.00	
79) Terphenyl-d14	19.503	244	1787654	83.703	ng	0.00	
<b>Target Compounds</b>							
2) 1,4-Dioxane	3.187	88	43939	40.796	ng		Qvalue 97
3) Pyridine	3.616	79	132966m	44.143	ng		
4) n-Nitrosodimethylamine	3.539	42	51359	43.058	ng		99
6) Aniline	7.000	93	99499	33.409	ng		97
8) 2-Chlorophenol	7.206	128	139458	40.672	ng		98
9) Benzaldehyde	6.765	77	67268	34.656	ng		99
10) Phenol	6.965	94	179912	41.653	ng		99
11) bis(2-Chloroethyl)ether	7.024	93	165037m	46.155	ng		
12) 1,3-Dichlorobenzene	7.447	146	149848	40.030	ng		99
13) 1,4-Dichlorobenzene	7.593	146	151916	40.067	ng		99
14) 1,2-Dichlorobenzene	7.922	146	149116	40.065	ng		99
15) Benzyl Alcohol	7.893	79	98580	42.529	ng		99
16) 2,2'-oxybis(1-Chloropr...	8.063	45	174241	41.166	ng		96
17) 2-Methylphenol	8.146	107	113951	40.749	ng		99
18) Hexachloroethane	8.616	117	50871	39.729	ng		98
19) n-Nitroso-di-n-propyla...	8.363	70	110591	42.268	ng		99
20) 3+4-Methylphenols	8.486	107	161794	41.727	ng		97
22) Acetophenone	8.410	105	215384	40.043	ng	#	99
24) Nitrobenzene	8.810	77	163457	41.795	ng		97
25) Isophorone	9.262	82	305817	41.792	ng		99
26) 2-Nitrophenol	9.497	139	85922	41.283	ng		99
27) 2,4-Dimethylphenol	9.615	122	96588	40.108	ng		97
28) bis(2-Chloroethoxy)met...	9.756	93	180210	40.232	ng		99
29) 2,4-Dichlorophenol	10.055	162	139985	40.947	ng		98
30) 1,2,4-Trichlorobenzene	10.173	180	148845	39.001	ng		100
31) Naphthalene	10.373	128	479224	39.967	ng		100
32) Benzoic acid	9.897	122	94456	42.834	ng		99
33) 4-Chloroaniline	10.602	127	178461	42.324	ng		98
34) Hexachlorobutadiene	10.625	225	90206	38.348	ng		98
35) Caprolactam	11.383	113	57396	45.719	ng		90
36) 4-Chloro-3-methylphenol	11.783	107	162513	43.532	ng		98
37) 2-Methylnaphthalene	11.965	142	349131	40.995	ng		98
38) 1-Methylnaphthalene	12.194	142	347270	40.889	ng		100
40) 1,2,4,5-Tetrachloroben...	12.323	216	175622	37.079	ng		100
41) Hexachlorocyclopentadiene	12.306	237	54568	39.479	ng		99
43) 2,4,6-Trichlorophenol	12.646	196	128512	39.507	ng		98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.752	196	144979	40.006	ng	97
46) 1,1'-Biphenyl	12.999	154	476785	38.475	ng	99
47) 2-Chloronaphthalene	13.046	162	376267	38.497	ng	99
48) 2-Nitroaniline	13.375	65	114738	44.348	ng	94
49) Acenaphthylene	13.904	152	588289	40.390	ng	99
50) Dimethylphthalate	13.675	163	519405	40.601	ng	100
51) 2,6-Dinitrotoluene	13.816	165	123306	41.761	ng	99
52) Acenaphthene	14.233	154	375195	40.367	ng	99
53) 3-Nitroaniline	14.227	138	127306	44.412	ng	100
54) 2,4-Dinitrophenol	14.362	184	79363	45.866	ng	96
55) Dibenzofuran	14.579	168	600773	40.139	ng	99
56) 4-Nitrophenol	14.650	139	106187	44.862	ng	93
57) 2,4-Dinitrotoluene	14.603	165	178718	43.083	ng	99
58) Fluorene	15.214	166	516260	41.330	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.850	232	136587	40.732	ng	99
60) Diethylphthalate	15.002	149	559909	41.509	ng	100
61) 4-Chlorophenyl-phenyle...	15.202	204	254962	40.196	ng	99
62) 4-Nitroaniline	15.414	138	140722	45.566	ng	97
63) Azobenzene	15.496	77	472139	42.904	ng	97
65) 4,6-Dinitro-2-methylph...	15.343	198	110950	44.402	ng	94
66) n-Nitrosodiphenylamine	15.455	169	451302	40.398	ng	100
67) 4-Bromophenyl-phenylether	16.089	248	180499	38.878	ng	98
68) Hexachlorobenzene	16.195	284	245701	40.211	ng	98
69) Atrazine	16.389	200	141676	42.956	ng	98
70) Pentachlorophenol	16.606	266	141175	42.583	ng	99
71) Phenanthrene	16.953	178	823366	40.216	ng	99
72) Anthracene	17.041	178	823243	40.718	ng	100
73) Carbazole	17.370	167	823854	40.633	ng	99
74) Di-n-butylphthalate	17.811	149	1011598	40.401	ng	100
75) Fluoranthene	18.957	202	1051631	40.062	ng	100
77) Benzidine	19.209	184	393870	38.785	ng	100
78) Pyrene	19.327	202	1109254	41.238	ng	100
80) Butylbenzylphthalate	20.173	149	492931	40.778	ng	98
81) Benzo(a)anthracene	21.054	228	1150750	40.986	ng	100
82) 3,3'-Dichlorobenzidine	21.025	252	449864	40.702	ng	99
83) Chrysene	21.107	228	1076464	40.336	ng	99
84) Bis(2-ethylhexyl)phtha...	20.872	149	721606	39.484	ng	100
85) Di-n-octyl phthalate	21.718	149	1227903	39.714	ng	99
87) Indeno(1,2,3-cd)pyrene	25.690	276	1665892	40.238	ng	99
88) Benzo(b)fluoranthene	22.652	252	1296785	39.233	ng	100
89) Benzo(k)fluoranthene	22.693	252	1255242	41.834	ng	100
90) Benzo(a)pyrene	23.252	252	1154661	40.463	ng	100
91) Dibenzo(a,h)anthracene	25.684	278	1396259	40.495	ng	100
92) Benzo(g,h,i)perylene	26.436	276	1379884	39.391	ng	99

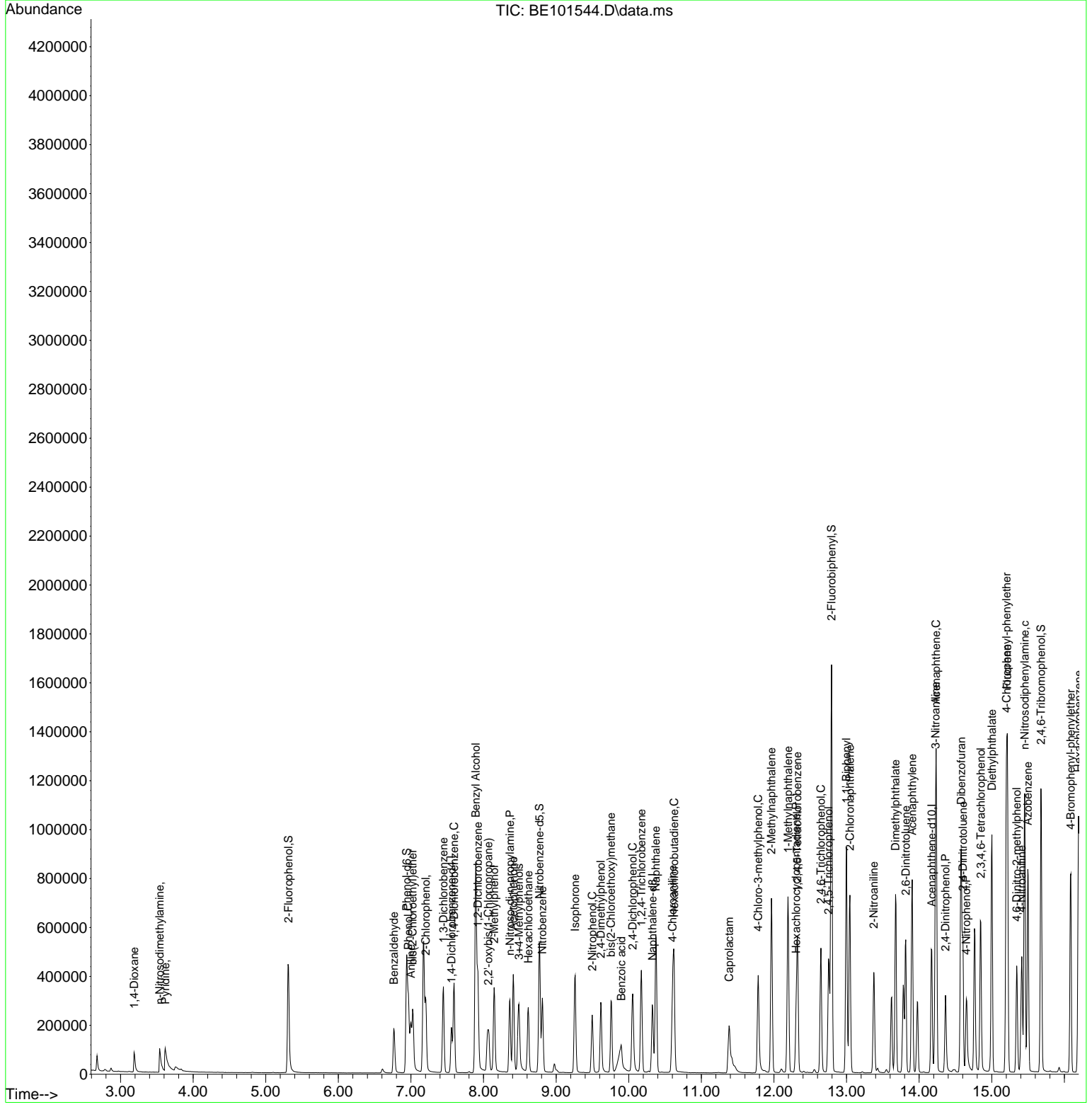
(#) = qualifier out of range (m) = manual integration (+) = signals summed

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