

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102324\
 Data File : BM048202.D
 Acq On : 23 Oct 2024 13:45
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 10/24/2024

Supervised By :mohammad ahmed 10/24/2024

Quant Time: Oct 23 17:58:10 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102324.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Oct 23 17:45:46 2024

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	164214	20.000	ng	0.00
21) Naphthalene-d8	10.516	136	675806	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	430191	20.000	ng	0.00
64) Phenanthrene-d10	17.110	188	872805	20.000	ng	0.00
76) Chrysene-d12	21.339	240	829636	20.000	ng	0.00
86) Perylene-d12	24.291	264	941862	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	194083	19.867	ng	0.00
7) Phenol-d6	6.904	99	253727	19.943	ng	0.00
23) Nitrobenzene-d5	8.881	82	235450	19.672	ng	0.00
42) 2,4,6-Tribromophenol	15.857	330	105824	20.097	ng	0.00
45) 2-Fluorobiphenyl	12.986	172	584048	20.833	ng	0.00
79) Terphenyl-d14	19.733	244	832028	20.466	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.205	88	43417	10.857	ng	99
3) Pyridine	3.616	79	108898	10.268	ng	98
4) n-Nitrosodimethylamine	3.516	42	45055	9.989	ng	# 95
6) Aniline	7.051	93	112524	10.655	ng	98
8) 2-Chlorophenol	7.293	128	105247	9.916	ng	99
9) Benzaldehyde	6.869	77	71283m	11.385	ng	
10) Phenol	6.934	94	126306	9.868	ng	97
11) bis(2-Chloroethyl)ether	7.151	93	104704	10.270	ng	96
12) 1,3-Dichlorobenzene	7.610	146	125778	10.279	ng	100
13) 1,4-Dichlorobenzene	7.757	146	128985	10.376	ng	98
14) 1,2-Dichlorobenzene	8.075	146	125469	10.444	ng	98
15) Benzyl Alcohol	7.969	79	77448	9.542	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.257	45	153954	10.324	ng	99
17) 2-Methylphenol	8.181	107	84745	9.963	ng	98
18) Hexachloroethane	8.798	117	46396	10.183	ng	97
19) n-Nitroso-di-n-propyla...	8.528	70	81492	10.125	ng	98
20) 3+4-Methylphenols	8.510	107	112836	9.576	ng	97
22) Acetophenone	8.545	105	165016	10.005	ng	# 97
24) Nitrobenzene	8.922	77	122626	9.888	ng	97
25) Isophorone	9.451	82	215452	9.820	ng	99
26) 2-Nitrophenol	9.634	139	50967	8.901	ng	94
27) 2,4-Dimethylphenol	9.704	122	65810	9.473	ng	96
28) bis(2-Chloroethoxy)met...	9.934	93	136287	9.817	ng	100
29) 2,4-Dichlorophenol	10.175	162	94106	9.417	ng	95
30) 1,2,4-Trichlorobenzene	10.381	180	113193	9.890	ng	97
31) Naphthalene	10.563	128	354273	10.051	ng	99
32) Benzoic acid	9.810	122	64616m	8.320	ng	
33) 4-Chloroaniline	10.686	127	111559	9.759	ng	99
34) Hexachlorobutadiene	10.851	225	71791	10.120	ng	99
35) Caprolactam	11.469	113	29958m	9.243	ng	
36) 4-Chloro-3-methylphenol	11.822	107	99967	9.582	ng	99
37) 2-Methylnaphthalene	12.180	142	234601	9.962	ng	99
38) 1-Methylnaphthalene	12.404	142	233676	10.007	ng	100
40) 1,2,4,5-Tetrachloroben...	12.557	216	123639	10.055	ng	99
41) Hexachlorocyclopentadiene	12.533	237	40205	9.663	ng	97
43) 2,4,6-Trichlorophenol	12.804	196	79475	9.740	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.880	196	87457	9.611	ng	98
46) 1,1'-Biphenyl	13.198	154	313591	10.210	ng	99
47) 2-Chloronaphthalene	13.239	162	244337	10.089	ng	98
48) 2-Nitroaniline	13.451	65	59910	9.121	ng	97
49) Acenaphthylene	14.092	152	351454	9.947	ng	99
50) Dimethylphthalate	13.827	163	297690	10.137	ng	99
51) 2,6-Dinitrotoluene	13.945	165	62280	9.504	ng	91
52) Acenaphthene	14.433	154	228492	10.178	ng	98
53) 3-Nitroaniline	14.280	138	56940	9.214	ng	# 94
54) 2,4-Dinitrophenol	14.492	184	29288	7.649	ng	99
55) Dibenzofuran	14.769	168	367889	10.315	ng	99
56) 4-Nitrophenol	14.610	139	43683	7.920	ng	95
57) 2,4-Dinitrotoluene	14.739	165	79851	9.228	ng	94
58) Fluorene	15.416	166	292703	10.378	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.004	232	75952	10.070	ng	99
60) Diethylphthalate	15.192	149	298826	10.050	ng	99
61) 4-Chlorophenyl-phenyle...	15.416	204	147671	10.472	ng	97
62) 4-Nitroaniline	15.445	138	53233	8.324	ng	95
63) Azobenzene	15.704	77	266197	9.989	ng	99
65) 4,6-Dinitro-2-methylph...	15.504	198	42770	8.446	ng	88
66) n-Nitrosodiphenylamine	15.627	169	242589	10.125	ng	99
67) 4-Bromophenyl-phenylether	16.304	248	89539	10.118	ng	96
68) Hexachlorobenzene	16.421	284	108457	10.205	ng	99
69) Atrazine	16.580	200	74922	11.175	ng	98
70) Pentachlorophenol	16.768	266	65591	9.400	ng	99
71) Phenanthrene	17.151	178	441552	10.217	ng	100
72) Anthracene	17.245	178	415460	9.930	ng	100
73) Carbazole	17.521	167	402900	9.824	ng	99
74) Di-n-butylphthalate	18.080	149	486204	9.524	ng	99
75) Fluoranthene	19.168	202	517157	10.031	ng	100
77) Benzidine	19.368	184	111510	11.646	ng	99
78) Pyrene	19.533	202	544382	9.803	ng	99
80) Butylbenzylphthalate	20.427	149	208443	8.977	ng	94
81) Benzo(a)anthracene	21.321	228	521872	9.844	ng	99
82) 3,3'-Dichlorobenzidine	21.245	252	180293	9.928	ng	99
83) Chrysene	21.380	228	507606	9.975	ng	98
84) Bis(2-ethylhexyl)phtha...	21.245	149	314375	9.315	ng	99
85) Di-n-octyl phthalate	22.356	149	506153	8.659	ng	99
87) Indeno(1,2,3-cd)pyrene	27.615	276	606458	9.737	ng	100
88) Benzo(b)fluoranthene	23.356	252	523956	9.702	ng	99
89) Benzo(k)fluoranthene	23.421	252	539599	10.335	ng	99
90) Benzo(a)pyrene	24.150	252	454217	9.675	ng	99
91) Dibenzo(a,h)anthracene	27.668	278	505969	9.755	ng	99
92) Benzo(g,h,i)perylene	28.650	276	517660	9.747	ng	98

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

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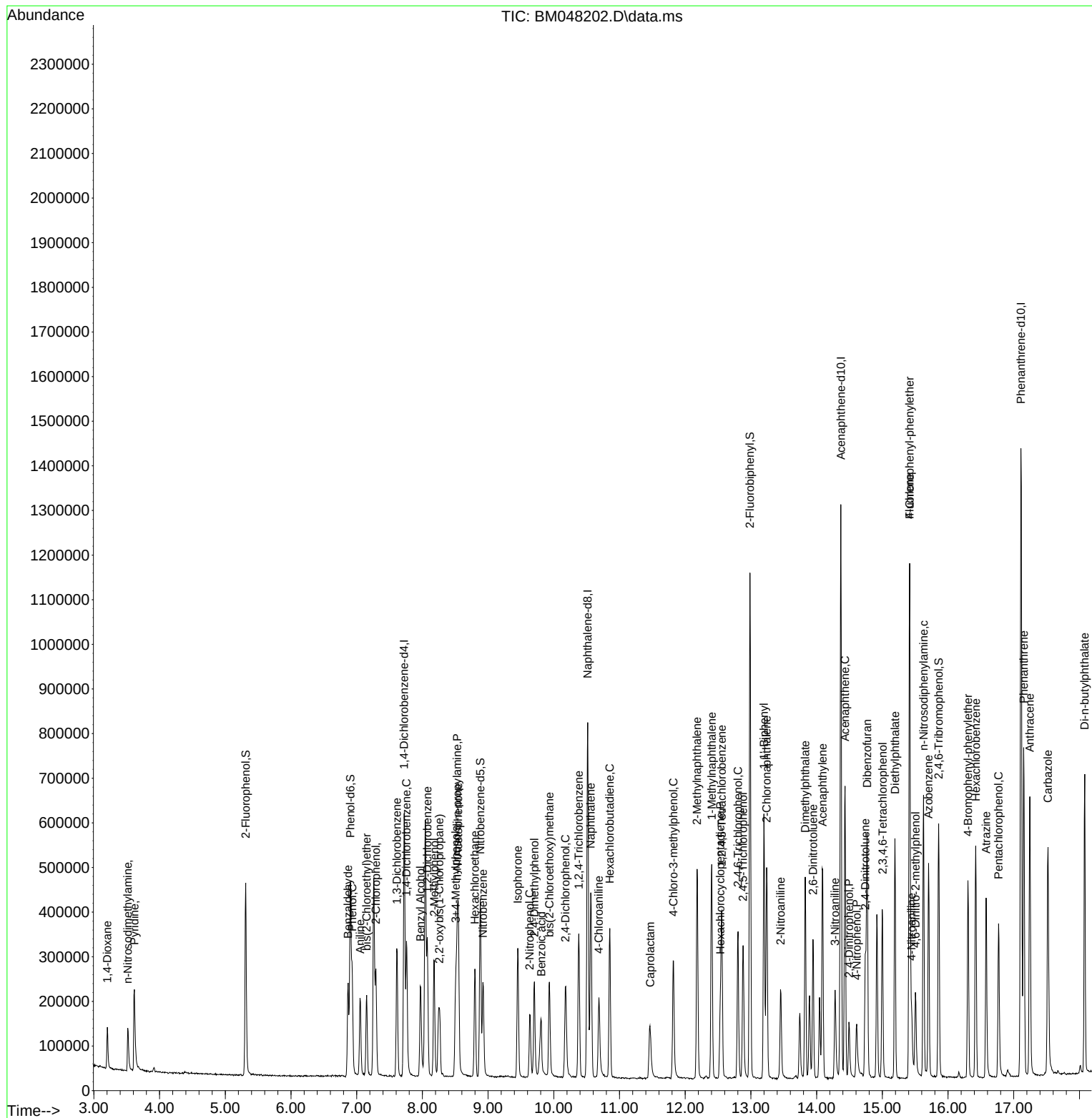
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