

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102324\
 Data File : BM048201.D
 Acq On : 23 Oct 2024 13:05
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDICC005

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 10/24/2024

Supervised By :mohammad ahmed 10/24/2024

Quant Time: Oct 23 17:57:13 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	145816	20.000	ng	0.00
21) Naphthalene-d8	10.516	136	593176	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	384226	20.000	ng	0.00
64) Phenanthrene-d10	17.110	188	791833	20.000	ng	0.00
76) Chrysene-d12	21.339	240	785602	20.000	ng	0.00
86) Perylene-d12	24.291	264	908010	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	90271	10.406	ng	0.00
7) Phenol-d6	6.904	99	113363	10.035	ng	0.00
23) Nitrobenzene-d5	8.881	82	102811	9.786	ng	0.00
42) 2,4,6-Tribromophenol	15.863	330	46245	9.833	ng	0.00
45) 2-Fluorobiphenyl	12.986	172	267303	10.675	ng	0.00
79) Terphenyl-d14	19.733	244	397731	10.332	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.205	88	19616	5.524	ng	98
3) Pyridine	3.616	79	47808	5.077	ng	# 89
4) n-Nitrosodimethylamine	3.522	42	20823	5.199	ng	# 97
6) Aniline	7.057	93	50475	5.383	ng	97
8) 2-Chlorophenol	7.293	128	48723	5.170	ng	97
9) Benzaldehyde	6.869	77	31542m	5.674	ng	
10) Phenol	6.934	94	56303	4.954	ng	97
11) bis(2-Chloroethyl)ether	7.151	93	45124	4.984	ng	99
12) 1,3-Dichlorobenzene	7.616	146	58340	5.369	ng	100
13) 1,4-Dichlorobenzene	7.763	146	59711	5.409	ng	97
14) 1,2-Dichlorobenzene	8.075	146	57354	5.376	ng	99
15) Benzyl Alcohol	7.975	79	32867	4.560	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.257	45	71883	5.429	ng	100
17) 2-Methylphenol	8.181	107	36152	4.787	ng	98
18) Hexachloroethane	8.798	117	21944	5.424	ng	95
19) n-Nitroso-di-n-propyla...	8.528	70	37763	5.284	ng	98
20) 3+4-Methylphenols	8.510	107	51475	4.920	ng	95
22) Acetophenone	8.551	105	73092	5.049	ng	# 91
24) Nitrobenzene	8.928	77	54443	5.002	ng	98
25) Isophorone	9.451	82	96675	5.020	ng	97
26) 2-Nitrophenol	9.640	139	21494	4.277	ng	# 82
27) 2,4-Dimethylphenol	9.704	122	28457	4.667	ng	97
28) bis(2-Chloroethoxy)met...	9.934	93	59074	4.848	ng	99
29) 2,4-Dichlorophenol	10.187	162	40064	4.567	ng	94
30) 1,2,4-Trichlorobenzene	10.381	180	51928	5.169	ng	98
31) Naphthalene	10.563	128	162694	5.259	ng	99
32) Benzoic acid	9.792	122	28049m	4.115	ng	
33) 4-Chloroaniline	10.692	127	47387	4.723	ng	99
34) Hexachlorobutadiene	10.851	225	32476	5.216	ng	96
35) Caprolactam	11.463	113	12948m	4.551	ng	
36) 4-Chloro-3-methylphenol	11.822	107	43099	4.707	ng	95
37) 2-Methylnaphthalene	12.186	142	107644	5.208	ng	99
38) 1-Methylnaphthalene	12.404	142	107297	5.235	ng	97
40) 1,2,4,5-Tetrachloroben...	12.557	216	56505	5.145	ng	99
41) Hexachlorocyclopentadiene	12.533	237	17150	4.615	ng	96
43) 2,4,6-Trichlorophenol	12.804	196	34246	4.699	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.886	196	39495	4.860	ng	97
46) 1,1'-Biphenyl	13.198	154	141745	5.167	ng	98
47) 2-Chloronaphthalene	13.239	162	110771	5.121	ng	99
48) 2-Nitroaniline	13.457	65	24776	4.223	ng	90
49) Acenaphthylene	14.092	152	159529	5.055	ng	98
50) Dimethylphthalate	13.827	163	139365	5.313	ng	99
51) 2,6-Dinitrotoluene	13.945	165	26245	4.484	ng	86
52) Acenaphthene	14.433	154	105474	5.260	ng	99
53) 3-Nitroaniline	14.286	138	22015	3.989	ng #	98
55) Dibenzofuran	14.769	168	169629	5.325	ng	100
57) 2,4-Dinitrotoluene	14.745	165	32425	4.195	ng	93
58) Fluorene	15.416	166	135671	5.386	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.004	232	34166	5.072	ng	99
60) Diethylphthalate	15.192	149	138791	5.226	ng	99
61) 4-Chlorophenyl-phenyle...	15.416	204	67194	5.335	ng	97
62) 4-Nitroaniline	15.451	138	20429	3.576	ng	94
63) Azobenzene	15.704	77	114836	4.825	ng	97
65) 4,6-Dinitro-2-methylph...	15.510	198	17234	3.751	ng #	80
66) n-Nitrosodiphenylamine	15.627	169	108646	4.998	ng	99
67) 4-Bromophenyl-phenylether	16.304	248	40550	5.051	ng	99
68) Hexachlorobenzene	16.421	284	49891	5.174	ng	98
69) Atrazine	16.580	200	34838	5.728	ng	98
70) Pentachlorophenol	16.774	266	28634	4.523	ng	98
71) Phenanthrene	17.151	178	206538	5.268	ng	99
72) Anthracene	17.245	178	190046	5.007	ng	99
73) Carbazole	17.527	167	180407	4.848	ng	96
74) Di-n-butylphthalate	18.080	149	223946	4.835	ng	98
75) Fluoranthene	19.174	202	244728	5.232	ng	99
77) Benzidine	19.368	184	81971	9.041	ng	98
78) Pyrene	19.533	202	262294	4.988	ng	99
80) Butylbenzylphthalate	20.433	149	96275	4.378	ng	100
81) Benzo(a)anthracene	21.321	228	255193	5.083	ng	99
82) 3,3'-Dichlorobenzidine	21.251	252	84774	4.930	ng	98
83) Chrysene	21.380	228	249978	5.188	ng	99
84) Bis(2-ethylhexyl)phtha...	21.245	149	144401	4.518	ng	98
85) Di-n-octyl phthalate	22.356	149	236872	4.280	ng #	99
87) Indeno(1,2,3-cd)pyrene	27.621	276	293568	4.889	ng	99
88) Benzo(b)fluoranthene	23.362	252	259343	4.981	ng	99
89) Benzo(k)fluoranthene	23.421	252	257436	5.114	ng	99
90) Benzo(a)pyrene	24.156	252	223092	4.929	ng	100
91) Dibenzo(a,h)anthracene	27.680	278	248009	4.960	ng	98
92) Benzo(g,h,i)perylene	28.662	276	256774	5.015	ng	96

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

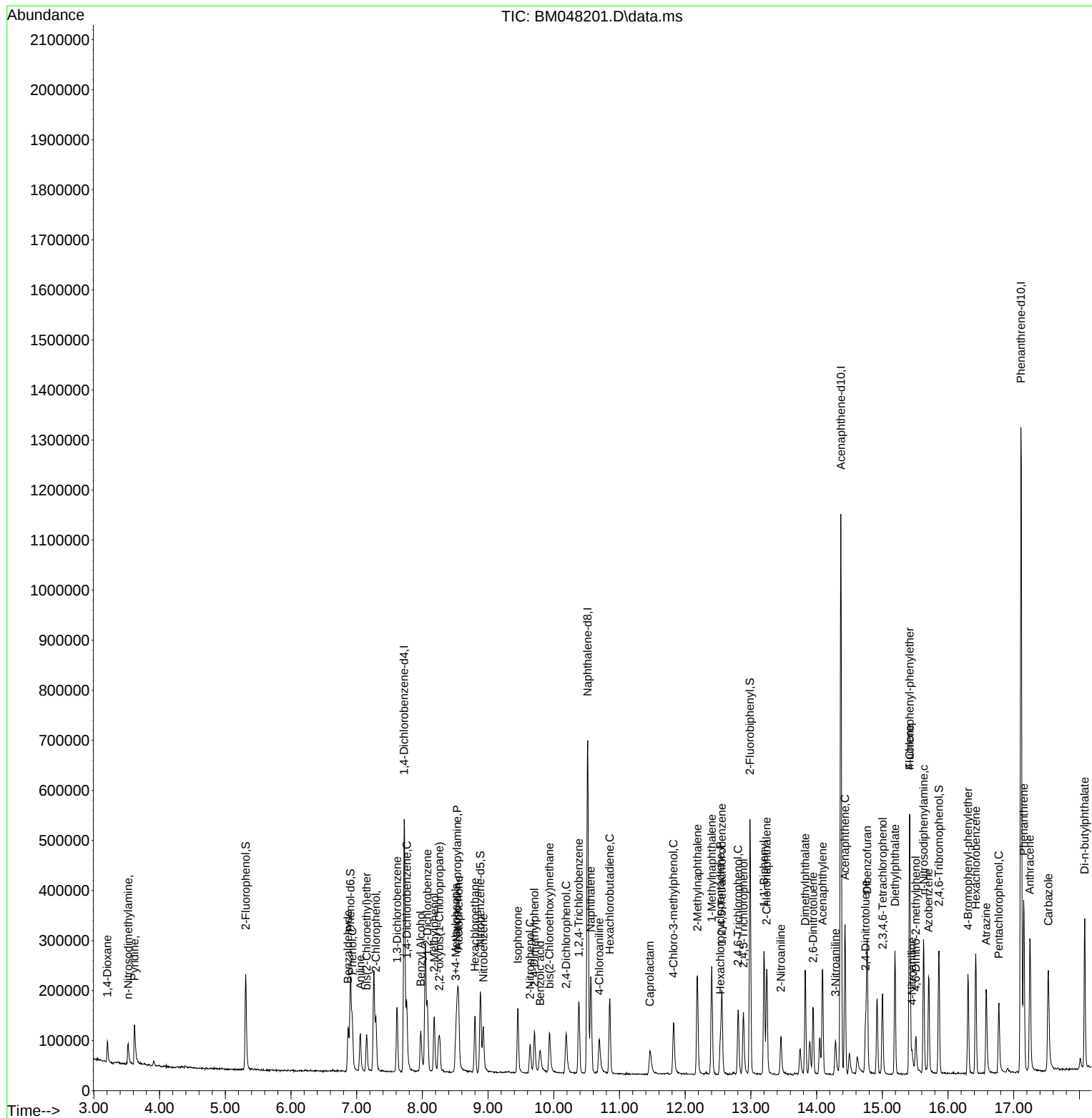
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