

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110424\
 Data File : BM048442.D
 Acq On : 05 Nov 2024 00:29
 Operator : RC/JU
 Sample : PB164575BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS575

Quant Time: Nov 05 01:00:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:17 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	3732	0.400	ng/ul	0.00
4) Naphthalene-d8	10.517	136	12853	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.357	164	6966	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.115	188	12683	0.400	ng/ul	0.00
17) Chrysene-d12	21.292	240	10685	0.400	ng/ul	0.00
23) Perylene-d12	23.531	264	10311	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	5027	1.097	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.134	152	7744	0.474	ng/ul	0.00
18) Fluoranthene-d10	19.140	212	15379	0.567	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.198	88	13530	2.721	ng/ul#	81
5) Naphthalene	10.567	128	15555	0.463	ng/ul	99
7) 2-Methylnaphthalene	12.206	142	8927	0.447	ng/ul	100
8) 1-Methylnaphthalene	12.404	142	11350	0.533	ng/ul	99
10) Acenaphthylene	14.080	152	12893	0.479	ng/ul	99
11) Acenaphthene	14.418	153	10299	0.477	ng/ul	98
12) Fluorene	15.426	166	11212	0.467	ng/ul	100
14) Pentachlorophenol	16.781	266	2930	0.830	ng/ul	94
15) Phenanthrene	17.157	178	15581	0.494	ng/ul	100
16) Anthracene	17.267	178	13843	0.469	ng/ul	99
19) Fluoranthene	19.168	202	20038	0.544	ng/ul	99
20) Pyrene	19.526	202	21486	0.543	ng/ul	98
21) Benzo(a)anthracene	21.284	228	14084	0.440	ng/ul	100
22) Chrysene	21.327	228	23977	0.544	ng/ul	100
24) Benzo(b)fluoranthene	22.856	252	17171	0.487	ng/ul	98
25) Benzo(k)fluoranthene	22.903	252	22379	0.536	ng/ul	99
26) Benzo(a)pyrene	23.440	252	16725	0.505	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.826	276	22058	0.477	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.840	278	17104	0.474	ng/ul	98
29) Benzo(g,h,i)perylene	26.500	276	19321	0.495	ng/ul	99

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

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