

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041324\
 Data File : BM045219.D
 Acq On : 13 Apr 2024 23:41
 Operator : MA/JU
 Sample : PB160169BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS169

Quant Time: Apr 15 00:37:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 15 00:32:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.631	152	1274	0.400	ng/ul	0.02
4) Naphthalene-d8	10.413	136	3815	0.400	ng/ul #	0.02
9) Acenaphthene-d10	14.270	164	1801	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.043	188	4120	0.400	ng/ul	0.01
17) Chrysene-d12	21.246	240	3471	0.400	ng/ul	0.00
23) Perylene-d12	23.467	264	4200	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	1154	0.690	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.024	152	1709	0.331	ng/ul	0.03
18) Fluoranthene-d10	19.075	212	3561	0.406	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.234	88	3079	1.919	ng/ul#	83
5) Naphthalene	10.462	128	3650	0.361	ng/ul	96
7) 2-Methylnaphthalene	12.101	142	2018	0.334	ng/ul	97
8) 1-Methylnaphthalene	12.305	142	2133	0.323	ng/ul	99
10) Acenaphthylene	14.001	152	3302	0.367	ng/ul	95
11) Acenaphthene	14.334	153	2440	0.379	ng/ul	97
12) Fluorene	15.352	166	2553	0.354	ng/ul	99
14) Pentachlorophenol	16.709	266	764	0.791	ng/ul	97
15) Phenanthrene	17.085	178	4100	0.353	ng/ul	98
16) Anthracene	17.191	178	4144	0.363	ng/ul	96
19) Fluoranthene	19.103	202	5141	0.414	ng/ul	96
20) Pyrene	19.470	202	5430	0.415	ng/ul	96
21) Benzo(a)anthracene	21.234	228	4320	0.358	ng/ul	96
22) Chrysene	21.281	228	6229	0.413	ng/ul	97
24) Benzo(b)fluoranthene	22.800	252	5407	0.356	ng/ul	96
25) Benzo(k)fluoranthene	22.847	252	7054	0.415	ng/ul	94
26) Benzo(a)pyrene	23.376	252	5691	0.393	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.716	276	7392	0.372	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.736	278	5824	0.370	ng/ul#	89
29) Benzo(g,h,i)perylene	26.376	276	6585	0.385	ng/ul	94

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

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