

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041324\
 Data File : BM045220.D
 Acq On : 14 Apr 2024 00:17
 Operator : MA/JU
 Sample : PB160163BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS163

Quant Time: Apr 15 00:37:28 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	7.635	152	1094	0.400	ng/ul	0.02
4) Naphthalene-d8	10.419	136	3086	0.400	ng/ul #	0.03
9) Acenaphthene-d10	14.274	164	1561	0.400	ng/ul	0.01
13) Phenanthrene-d10	17.047	188	3552	0.400	ng/ul	0.02
17) Chrysene-d12	21.249	240	3072	0.400	ng/ul	0.00
23) Perylene-d12	23.467	264	3707	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	1027	0.715	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.035	152	1506	0.360	ng/ul	0.04
18) Fluoranthene-d10	19.080	212	3399	0.438	ng/ul	0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.234	88	2940	2.134	ng/ul#	83
5) Naphthalene	10.468	128	3227	0.394	ng/ul#	93
7) 2-Methylnaphthalene	12.112	142	1823	0.373	ng/ul	98
8) 1-Methylnaphthalene	12.310	142	1914	0.358	ng/ul	97
10) Acenaphthylene	14.001	152	3051	0.392	ng/ul	96
11) Acenaphthene	14.339	153	2226	0.399	ng/ul	97
12) Fluorene	15.357	166	2418	0.387	ng/ul#	98
14) Pentachlorophenol	16.710	266	686	0.824	ng/ul	96
15) Phenanthrene	17.090	178	3841	0.384	ng/ul	95
16) Anthracene	17.199	178	3861	0.392	ng/ul	94
19) Fluoranthene	19.108	202	4895	0.446	ng/ul	97
20) Pyrene	19.471	202	5248	0.453	ng/ul	95
21) Benzo(a)anthracene	21.240	228	3963	0.371	ng/ul	98
22) Chrysene	21.287	228	6114	0.458	ng/ul	97
24) Benzo(b)fluoranthene	22.800	252	5144	0.384	ng/ul	94
25) Benzo(k)fluoranthene	22.847	252	6939	0.463	ng/ul	92
26) Benzo(a)pyrene	23.382	252	5296	0.414	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.719	276	6940	0.396	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.739	278	5347	0.385	ng/ul#	86
29) Benzo(g,h,i)perylene	26.377	276	6131	0.406	ng/ul	95

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

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