

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040824\
 Data File : BM045058.D
 Acq On : 09 Apr 2024 01:52
 Operator : MA/JU
 Sample : PB160056BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS056

Quant Time: Apr 09 03:34:30 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 : Tue Apr 09 02:13:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.639	152	1666	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.424	136	5745	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.284	164	3376	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.052	188	7533	0.400	ng/ul	0.00
17) Chrysene-d12	21.255	240	5797	0.400	ng/ul	0.00
23) Perylene-d12	23.473	264	6928	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	1165	0.532	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.030	152	2515	0.323	ng/ul	0.00
18) Fluoranthene-d10	19.085	212	5769	0.394	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.243	88	3229	1.539	ng/ul#	82
5) Naphthalene	10.474	128	5006	0.328	ng/ul	98
7) 2-Methylnaphthalene	12.112	142	3150	0.346	ng/ul	99
8) 1-Methylnaphthalene	12.316	142	3296	0.332	ng/ul	99
10) Acenaphthylene	14.011	152	5423	0.322	ng/ul	99
11) Acenaphthene	14.348	153	3959	0.328	ng/ul	98
12) Fluorene	15.357	166	4452	0.329	ng/ul	99
14) Pentachlorophenol	16.718	266	1234	0.699	ng/ul	94
15) Phenanthrene	17.094	178	7251	0.342	ng/ul	100
16) Anthracene	17.195	178	7120	0.341	ng/ul	98
19) Fluoranthene	19.117	202	8502	0.410	ng/ul	99
20) Pyrene	19.480	202	8796	0.402	ng/ul	97
21) Benzo(a)anthracene	21.240	228	7222	0.358	ng/ul	99
22) Chrysene	21.290	228	8547	0.339	ng/ul	99
24) Benzo(b)fluoranthene	22.809	252	9018	0.360	ng/ul	98
25) Benzo(k)fluoranthene	22.853	252	10210	0.364	ng/ul	96
26) Benzo(a)pyrene	23.385	252	8674	0.363	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.716	276	11678	0.357	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.743	278	9242	0.356	ng/ul	96
29) Benzo(g,h,i)perylene	26.390	276	10145	0.359	ng/ul	97

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

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