

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040824\
 Data File : BM045075.D
 Acq On : 09 Apr 2024 12:41
 Operator : MA/JU
 Sample : PB160067BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS067

Quant Time: Apr 09 13:10:58 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	7.631	152	1688	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.413	136	5642	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.279	164	3011	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.047	188	7012	0.400	ng/ul	0.00
17) Chrysene-d12	21.251	240	5725	0.400	ng/ul	0.00
23) Perylene-d12	23.469	264	7238	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	1236	0.557	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.024	152	2299	0.301	ng/ul	-0.01
18) Fluoranthene-d10	19.080	212	5560	0.384	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.243	88	3394	1.597	ng/ul#	83
5) Naphthalene	10.463	128	4769	0.319	ng/ul	97
7) 2-Methylnaphthalene	12.096	142	2999	0.336	ng/ul	99
8) 1-Methylnaphthalene	12.305	142	3069	0.314	ng/ul	99
10) Acenaphthylene	14.001	152	5040	0.335	ng/ul	99
11) Acenaphthene	14.344	153	3596	0.334	ng/ul	100
12) Fluorene	15.352	166	3960	0.328	ng/ul	99
14) Pentachlorophenol	16.705	266	1292	0.786	ng/ul	98
15) Phenanthrene	17.090	178	6264	0.317	ng/ul	98
16) Anthracene	17.187	178	6529	0.336	ng/ul	98
19) Fluoranthene	19.112	202	8115	0.397	ng/ul	99
20) Pyrene	19.475	202	8635	0.400	ng/ul	98
21) Benzo(a)anthracene	21.237	228	7213	0.362	ng/ul	98
22) Chrysene	21.286	228	8920	0.359	ng/ul	98
24) Benzo(b)fluoranthene	22.803	252	9056	0.346	ng/ul	98
25) Benzo(k)fluoranthene	22.847	252	10468	0.358	ng/ul	95
26) Benzo(a)pyrene	23.376	252	8933	0.358	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.705	276	12081	0.353	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.725	278	9482	0.350	ng/ul	96
29) Benzo(g,h,i)perylene	26.379	276	10383	0.352	ng/ul	97

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

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