

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070523\
 Data File : BM040709.D
 Acq On : 07 Jul 2023 11:56
 Operator : MA/JU
 Sample : PB153521BS
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS521

Quant Time: Jul 07 13:39:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 10:22:24 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	3444	0.400	ng/ul	0.00
4) Naphthalene-d8	10.708	136	14152	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.546	164	6982	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.303	188	13517	0.400	ng/ul	0.00
17) Chrysene-d12	21.492	240	8853	0.400	ng/ul	0.00
23) Perylene-d12	23.889	264	8460	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	2894	0.535	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.297	152	6459	0.326	ng/ul	0.00
18) Fluoranthene-d10	19.329	212	9958	0.351	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.339	88	7024	1.276	ng/ul	92
5) Naphthalene	10.757	128	12734	0.326	ng/ul	98
7) 2-Methylnaphthalene	12.374	142	7639	0.333	ng/ul	98
8) 1-Methylnaphthalene	12.588	142	7319	0.319	ng/ul	98
10) Acenaphthylene	14.268	152	9690	0.311	ng/ul	97
11) Acenaphthene	14.611	153	8223	0.312	ng/ul	98
12) Fluorene	15.601	166	8971	0.319	ng/ul	98
14) Pentachlorophenol	16.957	266	647	0.249	ng/ul	94
15) Phenanthrene	17.341	178	12799	0.305	ng/ul	98
16) Anthracene	17.434	178	10747	0.305	ng/ul	97
19) Fluoranthene	19.357	202	12968	0.345	ng/ul	99
20) Pyrene	19.724	202	13410	0.333	ng/ul	98
21) Benzo(a)anthracene	21.472	228	9618	0.321	ng/ul	99
22) Chrysene	21.524	228	10924	0.320	ng/ul	99
24) Benzo(b)fluoranthene	23.155	252	11029	0.337	ng/ul	93
25) Benzo(k)fluoranthene	23.202	252	10520	0.326	ng/ul#	90
26) Benzo(a)pyrene	23.784	252	9186	0.311	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.381	276	12745	0.310	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.394	278	9915	0.308	ng/ul	94
29) Benzo(g,h,i)perylene	27.142	276	11659	0.324	ng/ul	98

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

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