

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070823\
 Data File : BM040767.D
 Acq On : 09 Jul 2023 08:54
 Operator : MA/JU
 Sample : PB153891BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS891

Quant Time: Jul 09 14:31:16 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 : Sun Jul 09 14:07:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.891	152	3291	0.400	ng/ul	0.00
4) Naphthalene-d8	10.696	136	12372	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.536	164	5622	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.295	188	9827	0.400	ng/ul	0.00
17) Chrysene-d12	21.483	240	6169	0.400	ng/ul	0.00
23) Perylene-d12	23.880	264	5069	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	3521	0.681	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.291	152	5900	0.340	ng/ul	0.00
18) Fluoranthene-d10	19.324	212	7916	0.400	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	8808	1.674	ng/ul	90
5) Naphthalene	10.746	128	11737	0.344	ng/ul	98
7) 2-Methylnaphthalene	12.363	142	6867	0.343	ng/ul	98
8) 1-Methylnaphthalene	12.577	142	6420	0.321	ng/ul	98
10) Acenaphthylene	14.259	152	8597	0.342	ng/ul	96
11) Acenaphthene	14.601	153	7248	0.342	ng/ul	99
12) Fluorene	15.591	166	7709	0.340	ng/ul	97
14) Pentachlorophenol	16.953	266	587	0.310	ng/ul	96
15) Phenanthrene	17.337	178	10184	0.334	ng/ul	97
16) Anthracene	17.430	178	9024	0.352	ng/ul	94
19) Fluoranthene	19.352	202	10134	0.387	ng/ul	98
20) Pyrene	19.715	202	10712	0.381	ng/ul	97
21) Benzo(a)anthracene	21.469	228	6867	0.329	ng/ul	99
22) Chrysene	21.518	228	8216	0.345	ng/ul	98
24) Benzo(b)fluoranthene	23.146	252	7095	0.362	ng/ul	86
25) Benzo(k)fluoranthene	23.193	252	7513	0.389	ng/ul#	88
26) Benzo(a)pyrene	23.772	252	6247	0.354	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.364	276	8034	0.326	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.377	278	6291	0.326	ng/ul#	90
29) Benzo(g,h,i)perylene	27.122	276	7365	0.341	ng/ul	93

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

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