

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082423\
 Data File : BM041485.D
 Acq On : 25 Aug 2023 03:46
 Operator : MA/JU
 Sample : PB155032BS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS032

Quant Time: Aug 25 04:44:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM082423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 24 18:34:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.114	152	10075	0.400	ng/ul	0.00
4) Naphthalene-d8	10.944	136	26939	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.754	164	13099	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.506	188	26241	0.400	ng/ul	0.00
17) Chrysene-d12	21.696	240	9053	0.400	ng/ul	0.00
23) Perylene-d12	24.266	264	7659	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.448	96	8017	0.629	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.522	152	12256	0.344	ng/ul	0.00
18) Fluoranthene-d10	19.529	212	15649	0.392	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.486	88	18505	1.410	ng/ul#	91
5) Naphthalene	10.993	128	24562	0.306	ng/ul	100
7) 2-Methylnaphthalene	12.594	142	14839	0.308	ng/ul	100
8) 1-Methylnaphthalene	12.814	142	15884	0.324	ng/ul	99
10) Acenaphthylene	14.481	152	23203	0.295	ng/ul	100
11) Acenaphthene	14.819	153	16928	0.309	ng/ul	98
12) Fluorene	15.799	166	19155	0.313	ng/ul	100
14) Pentachlorophenol	17.126	266	5011	1.013	ng/ul	98
15) Phenanthrene	17.548	178	32774	0.340	ng/ul	99
16) Anthracene	17.641	178	35192	0.401	ng/ul	100
19) Fluoranthene	19.561	202	44136	0.531	ng/ul	99
20) Pyrene	19.924	202	46633	0.556	ng/ul	98
21) Benzo(a)anthracene	21.679	228	32926	0.559	ng/ul	99
22) Chrysene	21.734	228	36033	0.566	ng/ul	99
24) Benzo(b)fluoranthene	23.468	252	31717	0.605	ng/ul	98
25) Benzo(k)fluoranthene	23.523	252	30771	0.586	ng/ul	99
26) Benzo(a)pyrene	24.149	252	26959	0.626	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.984	276	29289	0.544	ng/ul#	98
28) Dibenzo(a,h)anthracene	27.018	278	21218	0.515	ng/ul	98
29) Benzo(g,h,i)perylene	27.829	276	24431	0.515	ng/ul	99

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

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