

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112423\
 Data File : BM043039.D
 Acq On : 26 Nov 2023 06:09
 Operator : MA/JU
 Sample : PB157100BS
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS100

Quant Time: Nov 27 02:46:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 27 02:36:19 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.872	152	569	0.400	ng/ul	0.00
4) Naphthalene-d8	10.640	136	2093	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	1210	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.230	188	2490	0.400	ng/ul	0.00
17) Chrysene-d12	21.417	240	1432	0.400	ng/ul	0.00
23) Perylene-d12	23.773	264	1833	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	954	1.093	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.251	152	1528	0.546	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	2657	0.509	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.197	88	2514	2.681	ng/ul#	83
5) Naphthalene	10.689	128	2605	0.458	ng/ul	98
7) 2-Methylnaphthalene	12.322	142	1574	0.469	ng/ul	100
8) 1-Methylnaphthalene	12.515	142	1597	0.427	ng/ul	100
10) Acenaphthylene	14.197	152	2841	0.460	ng/ul	97
11) Acenaphthene	14.525	153	2195	0.469	ng/ul	98
12) Fluorene	15.529	166	2284	0.475	ng/ul	98
14) Pentachlorophenol	16.892	266	610	0.905	ng/ul	96
15) Phenanthrene	17.276	178	3166	0.446	ng/ul	99
16) Anthracene	17.386	178	3046	0.452	ng/ul	98
19) Fluoranthene	19.290	202	3770	0.464	ng/ul	99
20) Pyrene	19.657	202	3988	0.462	ng/ul	98
21) Benzo(a)anthracene	21.406	228	2256	0.486	ng/ul	98
22) Chrysene	21.452	228	3966	0.458	ng/ul	99
24) Benzo(b)fluoranthene	23.042	252	3080	0.498	ng/ul	97
25) Benzo(k)fluoranthene	23.092	252	4247	0.499	ng/ul	97
26) Benzo(a)pyrene	23.674	252	3253	0.482	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.232	276	4471	0.489	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.255	278	3464	0.490	ng/ul	97
29) Benzo(g,h,i)perylene	26.989	276	4289	0.483	ng/ul	96

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

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