

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037644.D
 Acq On : 22 Nov 2022 16:23
 Operator : CG/JU
 Sample : PB149025BS
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS025

Quant Time: Nov 22 22:52:07 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM11622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 22:36:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.129	152	5206	0.400	ng/ul	0.00
4) Naphthalene-d8	10.947	136	14429	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.747	164	8461	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.479	188	19220	0.400	ng/ul	0.00
17) Chrysene-d12	21.639	240	13948	0.400	ng/ul	# 0.00
23) Perylene-d12	24.127	264	12793	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.450	96	4142	0.684	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.520	152	7325	0.306	ng/ul	0.00
18) Fluoranthene-d10	19.490	212	16331	0.354	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.488	88	10146	1.611	ng/ul#	51
5) Naphthalene	10.997	128	13832	0.316	ng/ul	99
7) 2-Methylnaphthalene	12.597	142	8827	0.296	ng/ul	99
8) 1-Methylnaphthalene	12.812	142	9517	0.312	ng/ul	99
10) Acenaphthylene	14.470	152	13170	0.280	ng/ul	100
11) Acenaphthene	14.812	153	10603	0.323	ng/ul	99
12) Fluorene	15.788	166	12451	0.315	ng/ul	99
14) Pentachlorophenol	17.124	266	2649	0.614	ng/ul	99
15) Phenanthrene	17.521	178	20979	0.319	ng/ul	100
16) Anthracene	17.614	178	18233	0.288	ng/ul	99
19) Fluoranthene	19.523	202	23813	0.354	ng/ul	96
20) Pyrene	19.885	202	24539	0.351	ng/ul	99
21) Benzo(a)anthracene	21.622	228	20534	0.329	ng/ul	99
22) Chrysene	21.677	228	22215	0.359	ng/ul	100
24) Benzo(b)fluoranthene	23.361	252	23551	0.404	ng/ul	90
25) Benzo(k)fluoranthene	23.411	252	21944	0.373	ng/ul#	90
26) Benzo(a)pyrene	24.013	252	19177	0.378	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.728	276	24052	0.358	ng/ul#	91
28) Dibenzo(a,h)anthracene	26.751	278	18510	0.351	ng/ul	94
29) Benzo(g,h,i)perylene	27.536	276	20301	0.356	ng/ul	96

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

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