

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037645.D
 Acq On : 22 Nov 2022 17:00
 Operator : CG/JU
 Sample : PB149043BS
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS043

Quant Time: Nov 22 22:52:17 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	5431	0.400	ng/ul	0.00
4) Naphthalene-d8	10.948	136	15084	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.747	164	8791	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.479	188	20274	0.400	ng/ul	0.00
17) Chrysene-d12	21.639	240	14489	0.400	ng/ul	# 0.00
23) Perylene-d12	24.124	264	13065	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.450	96	4290	0.679	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.526	152	7887	0.315	ng/ul	0.00
18) Fluoranthene-d10	19.490	212	17722	0.370	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.488	88	11162	1.698	ng/ul#	51
5) Naphthalene	10.997	128	15306	0.334	ng/ul	99
7) 2-Methylnaphthalene	12.597	142	9814	0.315	ng/ul	100
8) 1-Methylnaphthalene	12.812	142	10419	0.326	ng/ul	99
10) Acenaphthylene	14.470	152	14618	0.300	ng/ul	99
11) Acenaphthene	14.812	153	11752	0.345	ng/ul	98
12) Fluorene	15.788	166	13818	0.337	ng/ul	99
14) Pentachlorophenol	17.124	266	3024	0.665	ng/ul	99
15) Phenanthrene	17.521	178	23435	0.338	ng/ul	99
16) Anthracene	17.614	178	20152	0.301	ng/ul	99
19) Fluoranthene	19.523	202	26749	0.383	ng/ul	95
20) Pyrene	19.885	202	27319	0.376	ng/ul	99
21) Benzo(a)anthracene	21.622	228	22861	0.353	ng/ul	100
22) Chrysene	21.674	228	24793	0.386	ng/ul	99
24) Benzo(b)fluoranthene	23.361	252	26287	0.441	ng/ul	91
25) Benzo(k)fluoranthene	23.411	252	23858	0.397	ng/ul#	91
26) Benzo(a)pyrene	24.013	252	21024	0.406	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.731	276	26261	0.383	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.751	278	20355	0.377	ng/ul	95
29) Benzo(g,h,i)perylene	27.536	276	22282	0.382	ng/ul	96

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

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