

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037626.D
 Acq On : 22 Nov 2022 00:47
 Operator : CG/JU
 Sample : PB149187BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS187

Quant Time: Nov 22 01:37:20 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	5676	0.400	ng/ul	0.00
4) Naphthalene-d8	10.948	136	15699	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.747	164	9211	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.479	188	20776	0.400	ng/ul	0.00
17) Chrysene-d12	21.634	240	15424	0.400	ng/ul	# 0.00
23) Perylene-d12	24.121	264	14413	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.450	96	4618	0.700	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.520	152	8050	0.309	ng/ul	0.00
18) Fluoranthene-d10	19.490	212	17771	0.349	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.488	88	11523	1.678	ng/ul#	51
5) Naphthalene	10.997	128	16079	0.337	ng/ul	99
7) 2-Methylnaphthalene	12.597	142	10344	0.319	ng/ul	100
8) 1-Methylnaphthalene	12.812	142	11075	0.333	ng/ul	99
10) Acenaphthylene	14.470	152	15494	0.303	ng/ul	100
11) Acenaphthene	14.807	153	12470	0.349	ng/ul	98
12) Fluorene	15.788	166	14638	0.340	ng/ul	99
14) Pentachlorophenol	17.124	266	3370	0.723	ng/ul	98
15) Phenanthrene	17.521	178	24931	0.351	ng/ul	99
16) Anthracene	17.614	178	21794	0.318	ng/ul	100
19) Fluoranthene	19.523	202	29080	0.391	ng/ul	97
20) Pyrene	19.880	202	29832	0.386	ng/ul	97
21) Benzo(a)anthracene	21.619	228	25715	0.373	ng/ul	99
22) Chrysene	21.671	228	27941	0.408	ng/ul	100
24) Benzo(b)fluoranthene	23.355	252	30667	0.467	ng/ul	91
25) Benzo(k)fluoranthene	23.408	252	27683	0.417	ng/ul#	94
26) Benzo(a)pyrene	24.010	252	25409	0.445	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.725	276	31809	0.420	ng/ul	90
28) Dibenzo(a,h)anthracene	26.748	278	24485	0.412	ng/ul	95
29) Benzo(g,h,i)perylene	27.526	276	25860	0.402	ng/ul	96

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

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