

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035594.D
 Acq On : 22 Jun 2022 04:51
 Operator : CG/JU
 Sample : PB145592BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS592

Quant Time: Jun 22 05:38:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	3817	0.400	ng/ul	0.02
4) Naphthalene-d8	10.644	136	14875	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.469	164	7921	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	15526	0.400	ng/ul	0.00
17) Chrysene-d12	21.403	240	13258	0.400	ng/ul	0.00
23) Perylene-d12	23.748	264	12919	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.268	96	4515	0.841	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.244	152	8153	0.369	ng/ul	0.00
18) Fluoranthene-d10	19.247	212	16631	0.373	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.302	88	11521	2.155	ng/ul#	56
5) Naphthalene	10.694	128	15439	0.306	ng/ul	100
7) 2-Methylnaphthalene	12.321	142	9164	0.318	ng/ul	98
8) 1-Methylnaphthalene	12.519	142	9565	0.362	ng/ul	99
10) Acenaphthylene	14.196	152	14855	0.333	ng/ul	100
11) Acenaphthene	14.529	153	11040	0.334	ng/ul	99
12) Fluorene	15.533	166	12040	0.322	ng/ul	98
14) Pentachlorophenol	16.916	266	1920	0.645	ng/ul	100
15) Phenanthrene	17.267	178	18277	0.321	ng/ul	99
16) Anthracene	17.372	178	15754	0.316	ng/ul	99
19) Fluoranthene	19.275	202	22245	0.328	ng/ul	97
20) Pyrene	19.638	202	23825	0.335	ng/ul	97
21) Benzo(a)anthracene	21.392	228	16839	0.336	ng/ul	100
22) Chrysene	21.439	228	20552	0.345	ng/ul	99
24) Benzo(b)fluoranthene	23.034	252	19357	0.342	ng/ul	99
25) Benzo(k)fluoranthene	23.081	252	19689	0.335	ng/ul	99
26) Benzo(a)pyrene	23.648	252	20206	0.345	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.185	276	24224	0.359	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.192	278	19814	0.359	ng/ul	97
29) Benzo(g,h,i)perylene	26.920	276	21803	0.360	ng/ul	99

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

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