

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010322\
 Data File : BN018305.D
 Acq On : 03 Jan 2022 13:40
 Operator : CG/JU
 Sample : PB141605BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS605

Quant Time: Jan 04 00:00:12 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 28 00:34:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	1518	0.400	ng/ul	0.00
4) Naphthalene-d8	10.392	136	5387	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.253	164	3923	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.994	188	9535	0.400	ng/ul	0.00
17) Chrysene-d12	21.187	240	14661	0.400	ng/ul	0.00
23) Perylene-d12	23.408	264	17223	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.067	96	1043	0.736	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	2988	0.319	ng/ul	0.00
18) Fluoranthene-d10	19.026	212	13718	0.316	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.101	88	2723	1.727	ng/ul#	88
5) Naphthalene	10.442	128	5167	0.323	ng/ul	97
7) 2-Methylnaphthalene	12.064	142	3509	0.300	ng/ul	100
8) 1-Methylnaphthalene	12.284	142	3572	0.303	ng/ul	100
10) Acenaphthylene	13.970	152	5158	0.304	ng/ul	97
11) Acenaphthene	14.313	153	4595	0.336	ng/ul	98
12) Fluorene	15.303	166	5660	0.336	ng/ul	96
14) Pentachlorophenol	16.656	266	1916	0.508	ng/ul	100
15) Phenanthrene	17.032	178	10963	0.348	ng/ul	98
16) Anthracene	17.129	178	8853	0.312	ng/ul	97
19) Fluoranthene	19.054	202	17139	0.294	ng/ul	98
20) Pyrene	19.416	202	18814	0.309	ng/ul	98
21) Benzo(a)anthracene	21.169	228	18544	0.332	ng/ul	99
22) Chrysene	21.222	228	24016	0.396	ng/ul	99
24) Benzo(b)fluoranthene	22.739	252	25755	0.357	ng/ul	96
25) Benzo(k)fluoranthene	22.785	252	31235	0.380	ng/ul	97
26) Benzo(a)pyrene	23.312	252	25282	0.370	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.662	276	36135	0.415	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.682	278	27865	0.415	ng/ul	97
29) Benzo(g,h,i)perylene	26.349	276	33243	0.449	ng/ul	99

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

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