

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020222\
 Data File : BN018553.D
 Acq On : 02 Feb 2022 17:15
 Operator : CG/JU
 Sample : PB142349BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS349

Quant Time: Feb 03 03:31:53 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 14:08:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.927	152	379	0.400	ng/ul	0.00
4) Naphthalene-d8	10.735	136	1305	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.555	164	678	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.295	188	1297	0.400	ng/ul #	0.00
17) Chrysene-d12	21.466	240	1152	0.400	ng/ul #	0.00
23) Perylene-d12	23.860	264	982	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.295	96	300	0.689	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.319	152	633	0.356	ng/ul	0.00
18) Fluoranthene-d10	19.315	212	1324	0.402	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.333	88	806	1.713	ng/ul#	72
5) Naphthalene	10.784	128	1314	0.362	ng/ul#	90
7) 2-Methylnaphthalene	12.390	142	792	0.342	ng/ul	100
8) 1-Methylnaphthalene	12.610	142	810	0.341	ng/ul	99
10) Acenaphthylene	14.277	152	1046	0.388	ng/ul#	88
11) Acenaphthene	14.620	153	844	0.356	ng/ul	98
12) Fluorene	15.600	166	921	0.360	ng/ul#	96
14) Pentachlorophenol	16.945	266	134	0.554	ng/ul	97
15) Phenanthrene	17.337	178	1573	0.371	ng/ul#	92
16) Anthracene	17.426	178	1238	0.364	ng/ul#	89
19) Fluoranthene	19.343	202	1833	0.401	ng/ul#	83
20) Pyrene	19.706	202	1886	0.401	ng/ul#	85
21) Benzo(a)anthracene	21.451	228	1500	0.409	ng/ul	95
22) Chrysene	21.501	228	1943	0.433	ng/ul	97
24) Benzo(b)fluoranthene	23.129	252	1859	0.431	ng/ul	88
25) Benzo(k)fluoranthene	23.179	252	1847	0.409	ng/ul#	88
26) Benzo(a)pyrene	23.755	252	1601	0.460	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.344	276	1898	0.413	ng/ul#	79
28) Dibenzo(a,h)anthracene	26.368	278	1524	0.424	ng/ul#	78
29) Benzo(g,h,i)perylene	27.102	276	1766	0.431	ng/ul#	79

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

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