

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020122\
 Data File : BN018538.D
 Acq On : 02 Feb 2022 05:53
 Operator : CG/JU
 Sample : PB142390BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS390

Quant Time: Feb 02 06:22:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN013122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 03:59:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.927	152	417	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	1422	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.558	164	764	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.294	188	1511	0.400	ng/ul #	0.00
17) Chrysene-d12	21.465	240	1367	0.400	ng/ul #	0.00
23) Perylene-d12	23.861	264	1205	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	295	0.614	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	707	0.369	ng/ul	0.00
18) Fluoranthene-d10	19.319	212	1572	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.337	88	823	1.630	ng/ul#	75
5) Naphthalene	10.783	128	1364	0.348	ng/ul#	90
7) 2-Methylnaphthalene	12.394	142	847	0.341	ng/ul	99
8) 1-Methylnaphthalene	12.609	142	862	0.339	ng/ul	100
10) Acenaphthylene	14.276	152	1187	0.374	ng/ul#	89
11) Acenaphthene	14.618	153	941	0.356	ng/ul	98
12) Fluorene	15.603	166	1028	0.364	ng/ul#	97
14) Pentachlorophenol	16.947	266	232	1.728	ng/ul	100
15) Phenanthrene	17.336	178	1769	0.363	ng/ul#	93
16) Anthracene	17.424	178	1416	0.352	ng/ul#	91
19) Fluoranthene	19.346	202	2076	0.393	ng/ul#	86
20) Pyrene	19.709	202	2141	0.395	ng/ul#	87
21) Benzo(a)anthracene	21.450	228	1762	0.392	ng/ul	95
22) Chrysene	21.503	228	2150	0.415	ng/ul	97
24) Benzo(b)fluoranthene	23.130	252	2039	0.406	ng/ul	89
25) Benzo(k)fluoranthene	23.177	252	2093	0.397	ng/ul#	89
26) Benzo(a)pyrene	23.753	252	1850	0.443	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.349	276	2159	0.396	ng/ul#	80
28) Dibenzo(a,h)anthracene	26.370	278	1751	0.409	ng/ul#	80
29) Benzo(g,h,i)perylene	27.107	276	1973	0.416	ng/ul#	81

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

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