

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020222\
 Data File : BN018554.D
 Acq On : 02 Feb 2022 17:51
 Operator : CG/JU
 Sample : PB142390BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS390

Quant Time: Feb 03 03:32:07 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	312	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	1062	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.558	164	532	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.293	188	981	0.400	ng/ul #	0.00
17) Chrysene-d12	21.464	240	936	0.400	ng/ul #	0.00
23) Perylene-d12	23.861	264	817	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.294	96	283	0.790	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.317	152	602	0.416	ng/ul	0.00
18) Fluoranthene-d10	19.314	212	1226	0.458	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.332	88	791	2.042	ng/ul#	72
5) Naphthalene	10.783	128	1276	0.432	ng/ul#	89
7) 2-Methylnaphthalene	12.394	142	755	0.401	ng/ul	100
8) 1-Methylnaphthalene	12.608	142	771	0.398	ng/ul	99
10) Acenaphthylene	14.276	152	968	0.458	ng/ul#	87
11) Acenaphthene	14.618	153	794	0.427	ng/ul	98
12) Fluorene	15.603	166	844	0.420	ng/ul#	98
14) Pentachlorophenol	16.947	266	125	0.684	ng/ul	97
15) Phenanthrene	17.336	178	1420	0.442	ng/ul#	92
16) Anthracene	17.429	178	1097	0.426	ng/ul#	88
19) Fluoranthene	19.346	202	1681	0.452	ng/ul#	84
20) Pyrene	19.704	202	1735	0.454	ng/ul#	83
21) Benzo(a)anthracene	21.450	228	1395	0.468	ng/ul	94
22) Chrysene	21.502	228	1849	0.508	ng/ul	96
24) Benzo(b)fluoranthene	23.127	252	1797	0.500	ng/ul	88
25) Benzo(k)fluoranthene	23.177	252	1828	0.486	ng/ul#	87
26) Benzo(a)pyrene	23.753	252	1556	0.537	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.346	276	1905	0.498	ng/ul#	79
28) Dibenzo(a,h)anthracene	26.366	278	1538	0.514	ng/ul#	78
29) Benzo(g,h,i)perylene	27.104	276	1762	0.517	ng/ul#	79

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

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