

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012621\
 Data File : BN013486.D
 Acq On : 26 Jan 2021 13:06
 Operator : CG/JU
 Sample : PB134076BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS076

Quant Time: Jan 26 13:52:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN012521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 26 13:48:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.570	152	4117	0.40	ng/ul	0.00
4) Naphthalene-d8	10.329	136	14887	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.204	164	7661	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.949	188	16079	0.40	ng/ul	0.00
17) Chrysene-d12	21.154	240	14100	0.40	ng/ul	0.00
23) Perylene-d12	23.325	264	13600	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.178	96	3187	0.76	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.929	152	8060	0.37	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	16007	0.38	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.207	88	7969	1.90	ng/ul#	85
5) Naphthalene	10.378	128	15604	0.37	ng/ul	100
7) 2-Methylnaphthalene	12.001	142	10023	0.36	ng/ul	100
8) 1-Methylnaphthalene	12.226	142	9610	0.36	ng/ul	100
10) Acenaphthylene	13.917	152	11909	0.37	ng/ul	100
11) Acenaphthene	14.264	153	10349	0.37	ng/ul	99
12) Fluorene	15.259	166	11543	0.38	ng/ul	98
14) Pentachlorophenol	16.603	266	1485	0.76	ng/ul	99
15) Phenanthrene	16.992	178	19595	0.38	ng/ul	99
16) Anthracene	17.085	178	16322	0.39	ng/ul	99
19) Fluoranthene	19.018	202	21709	0.39	ng/ul	99
20) Pyrene	19.381	202	22018	0.39	ng/ul	98
21) Benzo(a)anthracene	21.139	228	18544	0.42	ng/ul	99
22) Chrysene	21.192	228	22051	0.40	ng/ul	100
24) Benzo(b)fluoranthene	22.679	252	21880	0.37	ng/ul	99
25) Benzo(k)fluoranthene	22.720	252	23799	0.38	ng/ul	98
26) Benzo(a)pyrene	23.232	252	19268	0.41	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.483	276	25241	0.38	ng/ul#	81
28) Dibenzo(a,h)anthracene	25.500	278	20310	0.38	ng/ul	100
29) Benzo(g,h,i)perylene	26.137	276	21539	0.37	ng/ul	98

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

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