

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012521\
 Data File : BN013481.D
 Acq On : 26 Jan 2021 01:30
 Operator : CG/JU
 Sample : SP5427
 Misc : SFAM SIM SPIKE
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP5427

Quant Time: Jan 26 02:37:31 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 : Mon Jan 25 22:48:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.574	152	5499	0.40	ng/ul	0.00
4) Naphthalene-d8	10.334	136	19585	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.204	164	10252	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	21964	0.40	ng/ul	0.00
17) Chrysene-d12	21.151	240	19477	0.40	ng/ul	0.00
23) Perylene-d12	23.323	264	15594	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.00	ng/ul	
6) 2-Methylnaphthalene-d10	0.000	152	0d	0.00	ng/ul	
18) Fluoranthene-d10	0.000	212	0d	0.00	ng/ul	
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.207	88	21153	3.77	ng/ul#	82
5) Naphthalene	10.378	128	37921	0.69	ng/ul	99
7) 2-Methylnaphthalene	12.001	142	24766	0.67	ng/ul	100
8) 1-Methylnaphthalene	12.226	142	23845	0.67	ng/ul	100
10) Acenaphthylene	13.917	152	27889	0.65	ng/ul	99
11) Acenaphthene	14.264	153	25021	0.68	ng/ul	100
12) Fluorene	15.254	166	28564	0.70	ng/ul	98
14) Pentachlorophenol	16.603	266	3835	1.45	ng/ul	99
15) Phenanthrene	16.992	178	47936	0.68	ng/ul	100
16) Anthracene	17.085	178	39686	0.69	ng/ul	99
19) Fluoranthene	19.019	202	48617	0.63	ng/ul	100
20) Pyrene	19.381	202	50617	0.65	ng/ul	99
21) Benzo(a)anthracene	21.137	228	41447	0.68	ng/ul	99
22) Chrysene	21.186	228	51130	0.67	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	47410	0.70	ng/ul	98
25) Benzo(k)fluoranthene	22.717	252	49001	0.69	ng/ul	97
26) Benzo(a)pyrene	23.226	252	36546	0.68	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.480	276	51808	0.69	ng/ul#	82
28) Dibenzo(a,h)anthracene	25.493	278	42759	0.70	ng/ul	98
29) Benzo(g,h,i)perylene	26.134	276	45554	0.68	ng/ul	98

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

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