

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062121\
 Data File : BN014952.D
 Acq On : 21 Jun 2021 12:49
 Operator : CG/JU
 Sample : SST0.801
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8001

Quant Time: Jun 21 13:22:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 21 13:17:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.747	152	3115	0.400	ng/ul	0.00
4) Naphthalene-d8	10.516	136	10589	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.366	164	5410	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.114	188	10678	0.400	ng/ul	0.00
17) Chrysene-d12	21.318	240	8580	0.400	ng/ul	0.00
23) Perylene-d12	23.574	264	6496	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	2778	0.627	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.111	152	13363	0.749	ng/ul	0.00
18) Fluoranthene-d10	19.149	212	23704	0.878	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	3053	0.743	ng/ul	91
5) Naphthalene	10.565	128	26699	0.795	ng/ul	99
7) 2-Methylnaphthalene	12.182	142	17580	0.796	ng/ul	99
8) 1-Methylnaphthalene	12.402	142	17264	0.809	ng/ul	98
10) Acenaphthylene	14.088	152	22060	0.774	ng/ul	100
11) Acenaphthene	14.431	153	16991	0.841	ng/ul	99
12) Fluorene	15.421	166	18960	0.819	ng/ul	99
14) Pentachlorophenol	16.764	266	1725	0.645	ng/ul	99
15) Phenanthrene	17.156	178	30798	0.865	ng/ul	100
16) Anthracene	17.249	178	26664	0.769	ng/ul	99
19) Fluoranthene	19.181	202	32745	0.953	ng/ul	99
20) Pyrene	19.544	202	32221	0.915	ng/ul	99
21) Benzo(a)anthracene	21.303	228	25401	0.818	ng/ul	100
22) Chrysene	21.353	228	30152	0.892	ng/ul	100
24) Benzo(b)fluoranthene	22.899	252	24578	0.965	ng/ul	99
25) Benzo(k)fluoranthene	22.945	252	27830	0.942	ng/ul	99
26) Benzo(a)pyrene	23.478	252	20428	0.876	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.852	276	24969	0.928	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.872	278	19665	0.919	ng/ul	99
29) Benzo(g,h,i)perylene	26.540	276	23245	0.942	ng/ul	100

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

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