

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090121\
 Data File : BN016120.D
 Acq On : 01 Sep 2021 08:26
 Operator : CG/JU
 Sample : SST00.115
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.1215

Quant Time: Sep 01 10:20:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 01 10:18:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	3108	0.400	ng/ul	0.00
4) Naphthalene-d8	10.655	136	11999	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	6572	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.247	188	13596	0.400	ng/ul	0.00
17) Chrysene-d12	21.444	240	13675	0.400	ng/ul	0.00
23) Perylene-d12	23.838	264	12914	0.400	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	377	0.089	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	1711	0.093	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	3786	0.104	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	539	0.123	ng/ul#	79
5) Naphthalene	10.710	128	3334	0.109	ng/ul	96
7) 2-Methylnaphthalene	12.321	142	2122	0.102	ng/ul	99
8) 1-Methylnaphthalene	12.541	142	2130	0.104	ng/ul	99
10) Acenaphthylene	14.220	152	2562	0.106	ng/ul	96
11) Acenaphthene	14.562	153	2182	0.111	ng/ul	99
12) Fluorene	15.548	166	2421	0.105	ng/ul	99
14) Pentachlorophenol	16.901	266	188	0.051	ng/ul	96
15) Phenanthrene	17.289	178	4126	0.110	ng/ul	98
16) Anthracene	17.378	178	3420	0.104	ng/ul	97
19) Fluoranthene	19.309	202	4588	0.110	ng/ul	96
20) Pyrene	19.671	202	4798	0.113	ng/ul	98
21) Benzo(a)anthracene	21.429	228	4370	0.101	ng/ul	97
22) Chrysene	21.482	228	5041	0.112	ng/ul	99
24) Benzo(b)fluoranthene	23.107	252	4994	0.108	ng/ul	87
25) Benzo(k)fluoranthene	23.157	252	5014	0.109	ng/ul#	86
26) Benzo(a)pyrene	23.730	252	4291	0.106	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.309	276	5463	0.103	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.329	278	4432	0.100	ng/ul#	90
29) Benzo(g,h,i)perylene	27.070	276	4776	0.106	ng/ul	95

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

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