

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015462.D
 Acq On : 10 Jul 2021 01:27
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4115

Quant Time: Jul 10 02:51:33 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 : Sat Jul 10 00:29:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	3262	0.400	ng/ul	0.00
4) Naphthalene-d8	10.479	136	11185	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.335	164	5702	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	11084	0.400	ng/ul	0.00
17) Chrysene-d12	21.281	240	8656	0.400	ng/ul	0.00
23) Perylene-d12	23.517	264	8057	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	1324	0.360	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.069	152	6805	0.388	ng/ul	0.00
18) Fluoranthene-d10	19.118	212	11438	0.402	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.315	88	1321	0.332	ng/ul	96
5) Naphthalene	10.529	128	14145	0.396	ng/ul	100
7) 2-Methylnaphthalene	12.146	142	9233	0.395	ng/ul	99
8) 1-Methylnaphthalene	12.366	142	8955	0.397	ng/ul	99
10) Acenaphthylene	14.053	152	11639	0.395	ng/ul	99
11) Acenaphthene	14.400	153	8941	0.396	ng/ul	99
12) Fluorene	15.390	166	9667	0.387	ng/ul	100
14) Pentachlorophenol	16.740	266	999	0.407	ng/ul	99
15) Phenanthrene	17.124	178	15904	0.395	ng/ul	99
16) Anthracene	17.217	178	13411	0.378	ng/ul	100
19) Fluoranthene	19.146	202	16288	0.407	ng/ul	98
20) Pyrene	19.513	202	16237	0.405	ng/ul	99
21) Benzo(a)anthracene	21.264	228	13365	0.417	ng/ul	99
22) Chrysene	21.316	228	15088	0.393	ng/ul	100
24) Benzo(b)fluoranthene	22.848	252	15331	0.407	ng/ul	94
25) Benzo(k)fluoranthene	22.889	252	15809	0.372	ng/ul	99
26) Benzo(a)pyrene	23.421	252	13872	0.439	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.767	276	17045	0.446	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.777	278	13433	0.449	ng/ul	97
29) Benzo(g,h,i)perylene	26.451	276	14145	0.400	ng/ul	99

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

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