

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071221\
 Data File : BN015525.D
 Acq On : 13 Jul 2021 13:02
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4121

Quant Time: Jul 13 13:33:00 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 Jul 12 11:02:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.808	152	3147	0.400	ng/ul	0.00
4) Naphthalene-d8	10.595	136	11085	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.441	164	5828	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.191	188	11418	0.400	ng/ul	0.00
17) Chrysene-d12	21.387	240	10594	0.400	ng/ul	0.00
23) Perylene-d12	23.742	264	9852	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	1262	0.356	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	6261	0.360	ng/ul	0.00
18) Fluoranthene-d10	19.225	212	11928	0.343	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	1322	0.345	ng/ul	97
5) Naphthalene	10.644	128	12731	0.360	ng/ul	100
7) 2-Methylnaphthalene	12.261	142	8298	0.359	ng/ul	99
8) 1-Methylnaphthalene	12.481	142	8140	0.364	ng/ul	99
10) Acenaphthylene	14.159	152	11575	0.384	ng/ul	99
11) Acenaphthene	14.506	153	8300	0.359	ng/ul	99
12) Fluorene	15.492	166	8880	0.348	ng/ul	97
14) Pentachlorophenol	16.845	266	1324	0.524	ng/ul	99
15) Phenanthrene	17.234	178	14944	0.361	ng/ul	100
16) Anthracene	17.322	178	13605	0.372	ng/ul	100
19) Fluoranthene	19.253	202	16976	0.347	ng/ul	96
20) Pyrene	19.615	202	17622	0.359	ng/ul	97
21) Benzo(a)anthracene	21.372	228	15598	0.397	ng/ul	96
22) Chrysene	21.422	228	15363	0.327	ng/ul	99
24) Benzo(b)fluoranthene	23.026	252	15112	0.328	ng/ul	96
25) Benzo(k)fluoranthene	23.073	252	15379	0.296	ng/ul	98
26) Benzo(a)pyrene	23.637	252	13443	0.348	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.162	276	16667	0.357	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.183	278	13151	0.359	ng/ul	98
29) Benzo(g,h,i)perylene	26.907	276	14044	0.325	ng/ul	99

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

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