

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016165.D
 Acq On : 03 Sep 2021 08:37
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4266

Quant Time: Sep 03 09:08:25 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 : Thu Sep 02 13:32:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	2773	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	11360	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	6481	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.250	188	12897	0.400	ng/ul	0.00
17) Chrysene-d12	21.439	240	13978	0.400	ng/ul	0.00
23) Perylene-d12	23.833	264	14455	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	1271	0.394	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	6624	0.414	ng/ul	0.00
18) Fluoranthene-d10	19.280	212	14822	0.395	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	1290	0.366	ng/ul#	92
5) Naphthalene	10.710	128	11525	0.381	ng/ul	100
7) 2-Methylnaphthalene	12.327	142	7849	0.398	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	7792	0.396	ng/ul	99
10) Acenaphthylene	14.224	152	10908	0.445	ng/ul	100
11) Acenaphthene	14.566	153	7947	0.387	ng/ul	99
12) Fluorene	15.552	166	8970	0.382	ng/ul	100
14) Pentachlorophenol	16.904	266	1186	0.568	ng/ul	98
15) Phenanthrene	17.293	178	14282	0.384	ng/ul	99
16) Anthracene	17.381	178	13837	0.431	ng/ul	99
19) Fluoranthene	19.313	202	17309	0.372	ng/ul	98
20) Pyrene	19.675	202	18089	0.384	ng/ul#	94
21) Benzo(a)anthracene	21.424	228	18806	0.440	ng/ul	99
22) Chrysene	21.477	228	18766	0.389	ng/ul	99
24) Benzo(b)fluoranthene	23.102	252	20328	0.374	ng/ul	96
25) Benzo(k)fluoranthene	23.152	252	19557	0.357	ng/ul	98
26) Benzo(a)pyrene	23.725	252	18389	0.399	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	26.310	276	23049	0.387	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.330	278	19104	0.389	ng/ul	98
29) Benzo(g,h,i)perylene	27.071	276	19410	0.375	ng/ul	99

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

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