

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090324\
 Data File : BN033747.D
 Acq On : 03 Sep 2024 19:22
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4303

Quant Time: Sep 04 01:42:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 01:39:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.519	152	6620	0.400	ng/ul	0.00
4) Naphthalene-d8	10.281	136	17824	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.159	164	8755	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.913	188	17823	0.400	ng/ul	0.00
17) Chrysene-d12	21.120	240	14385	0.400	ng/ul	0.00
23) Perylene-d12	23.273	264	15495	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.135	96	2767	0.407	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.876	152	9861	0.386	ng/ul	0.00
18) Fluoranthene-d10	18.951	212	17063	0.398	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.169	88	3098	0.410	ng/ul#	86
5) Naphthalene	10.331	128	18572	0.390	ng/ul	100
7) 2-Methylnaphthalene	11.953	142	11647	0.386	ng/ul	99
8) 1-Methylnaphthalene	12.179	142	11943	0.384	ng/ul	99
10) Acenaphthylene	13.877	152	15275	0.384	ng/ul	100
11) Acenaphthene	14.224	153	11603	0.404	ng/ul	99
12) Fluorene	15.219	166	12837	0.400	ng/ul	99
14) Pentachlorophenol	16.567	266	1612	0.287	ng/ul	99
15) Phenanthrene	16.956	178	19775	0.376	ng/ul	99
16) Anthracene	17.044	178	17602	0.372	ng/ul	99
19) Fluoranthene	18.979	202	21782	0.384	ng/ul	99
20) Pyrene	19.346	202	22785	0.378	ng/ul	98
21) Benzo(a)anthracene	21.105	228	21278	0.376	ng/ul	99
22) Chrysene	21.155	228	21775	0.381	ng/ul	99
24) Benzo(b)fluoranthene	22.633	252	22025	0.366	ng/ul	98
25) Benzo(k)fluoranthene	22.677	252	22935	0.381	ng/ul	98
26) Benzo(a)pyrene	23.180	252	18399	0.381	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.410	276	27281	0.376	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.420	278	21283	0.378	ng/ul	99
29) Benzo(g,h,i)perylene	26.057	276	21206	0.374	ng/ul	99

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

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