

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103124\
 Data File : BN034783.D
 Acq On : 01 Nov 2024 02:06
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4322

Quant Time: Nov 03 23:24:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 11:13:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.600	152	8571	0.400	ng/ul	0.00
4) Naphthalene-d8	10.366	136	26236	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.236	164	14401	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.979	188	29052	0.400	ng/ul	0.00
17) Chrysene-d12	21.172	240	18413	0.400	ng/ul	0.00
23) Perylene-d12	23.355	264	14209	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.174	96	3638	0.410	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.961	152	13332	0.405	ng/ul	0.00
18) Fluoranthene-d10	19.014	212	24358	0.406	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.208	88	4043	0.415	ng/ul#	88
5) Naphthalene	10.416	128	28210	0.406	ng/ul	100
7) 2-Methylnaphthalene	12.038	142	18443	0.408	ng/ul	99
8) 1-Methylnaphthalene	12.258	142	18712	0.407	ng/ul	99
10) Acenaphthylene	13.949	152	25994	0.397	ng/ul	99
11) Acenaphthene	14.296	153	20247	0.404	ng/ul	99
12) Fluorene	15.286	166	22396	0.401	ng/ul	100
14) Pentachlorophenol	16.633	266	1782	0.311	ng/ul	99
15) Phenanthrene	17.021	178	35869	0.408	ng/ul	100
16) Anthracene	17.110	178	32278	0.397	ng/ul	100
19) Fluoranthene	19.042	202	36828	0.412	ng/ul	99
20) Pyrene	19.404	202	37890	0.407	ng/ul	99
21) Benzo(a)anthracene	21.157	228	29075	0.395	ng/ul	100
22) Chrysene	21.210	228	29929	0.406	ng/ul	99
24) Benzo(b)fluoranthene	22.703	252	24735	0.428	ng/ul	100
25) Benzo(k)fluoranthene	22.747	252	24456	0.419	ng/ul	100
26) Benzo(a)pyrene	23.258	252	20231	0.406	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.527	276	22104	0.417	ng/ul	99
28) Dibenzo(a,h)anthracene	25.544	278	17049	0.425	ng/ul	99
29) Benzo(g,h,i)perylene	26.188	276	17606	0.423	ng/ul	99

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

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