

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034945.D
 Acq On : 09 Nov 2024 17:37
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4327

Quant Time: Nov 09 22:53:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.571	152	2450	0.400	ng/u1	0.00
4) Naphthalene-d8	10.333	136	6113	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.203	164	4775	0.400	ng/u1	0.00
13) Phenanthrene-d10	16.950	188	11993	0.400	ng/u1	0.00
17) Chrysene-d12	21.145	240	12317	0.400	ng/u1	0.00
23) Perylene-d12	23.311	264	12200	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	909	0.390	ng/u1	0.00
6) 2-Methylnaphthalene-d10	11.928	152	3842	0.410	ng/u1	0.00
18) Fluoranthene-d10	18.981	212	12810	0.405	ng/u1	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.182	88	964	0.377	ng/u1	97
5) Naphthalene	10.383	128	6094	0.398	ng/u1	98
7) 2-Methylnaphthalene	12.005	142	4623	0.405	ng/u1	100
8) 1-Methylnaphthalene	12.225	142	4656	0.406	ng/u1	98
10) Acenaphthylene	13.921	152	7488	0.384	ng/u1	100
11) Acenaphthene	14.268	153	5788	0.391	ng/u1	98
12) Fluorene	15.258	166	7233	0.384	ng/u1	99
14) Pentachlorophenol	16.603	266	1023	0.369	ng/u1	98
15) Phenanthrene	16.992	178	12638	0.394	ng/u1	99
16) Anthracene	17.080	178	11675	0.388	ng/u1	99
19) Fluoranthene	19.014	202	16395	0.398	ng/u1	99
20) Pyrene	19.376	202	17025	0.399	ng/u1	99
21) Benzo(a)anthracene	21.128	228	17141	0.384	ng/u1	99
22) Chrysene	21.180	228	17700	0.389	ng/u1	99
24) Benzo(b)fluoranthene	22.665	252	16581	0.387	ng/u1	100
25) Benzo(k)fluoranthene	22.709	252	17870	0.398	ng/u1	99
26) Benzo(a)pyrene	23.214	252	14699	0.384	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	25.463	276	17812	0.380	ng/u1	97
28) Dibenzo(a,h)anthracene	25.480	278	13931	0.381	ng/u1	98
29) Benzo(g,h,i)perylene	26.117	276	13446	0.376	ng/u1	98

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

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