

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034917.D
 Acq On : 08 Nov 2024 23:04
 Operator : RC/JU
 Sample : SSTD1.635
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6220

Quant Time: Nov 08 23:56:39 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 : Fri Nov 08 23:54:11 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	1805	0.400	ng/ul	0.00
4) Naphthalene-d8	10.339	136	4472	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.208	164	3468	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	8758	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	9667	0.400	ng/ul	0.00
23) Perylene-d12	23.314	264	9853	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	2644	1.415	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.934	152	10836	1.932	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	37667	1.197	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.187	88	2899	1.413	ng/ul#	94
5) Naphthalene	10.388	128	17331	1.463	ng/ul	96
7) 2-Methylnaphthalene	12.011	142	13176	1.711	ng/ul	99
8) 1-Methylnaphthalene	12.231	142	13293	1.695	ng/ul	98
10) Acenaphthylene	13.926	152	21765	1.379	ng/ul	98
11) Acenaphthene	14.268	153	16585	1.373	ng/ul	99
12) Fluorene	15.263	166	21326	1.585	ng/ul	98
14) Pentachlorophenol	16.607	266	3530	2.042	ng/ul	98
15) Phenanthrene	16.996	178	36781	1.389	ng/ul	98
16) Anthracene	17.084	178	35218	1.437	ng/ul	97
19) Fluoranthene	19.018	202	49709	1.060	ng/ul	100
20) Pyrene	19.381	202	51205	1.049	ng/ul	99
21) Benzo(a)anthracene	21.131	228	54674	1.413	ng/ul	99
22) Chrysene	21.183	228	54642	1.410	ng/ul	100
24) Benzo(b)fluoranthene	22.668	252	54524	1.360	ng/ul	95
25) Benzo(k)fluoranthene	22.711	252	56393	1.394	ng/ul#	94
26) Benzo(a)pyrene	23.220	252	48427	1.403	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.470	276	60609	1.649	ng/ul	98
28) Dibenzo(a,h)anthracene	25.487	278	47528	1.710	ng/ul	96
29) Benzo(g,h,i)perylene	26.124	276	45975	1.592	ng/ul	98

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

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