

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121724\
 Data File : BM049026.D
 Acq On : 17 Dec 2024 18:14
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4112

Quant Time: Dec 18 01:18:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM120924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 17 10:03:27 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	7945	0.400	ng/ul	0.00
4) Naphthalene-d8	10.337	136	22850	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.206	164	12824	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.951	188	24902	0.400	ng/ul #	0.00
17) Chrysene-d12	21.146	240	14857	0.400	ng/ul #	0.00
23) Perylene-d12	23.305	264	14367	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.155	96	3607	0.388	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.932	152	12099	0.415	ng/ul	0.00
18) Fluoranthene-d10	18.983	212	20803	0.469	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.189	88	4126	0.384	ng/ul	93
5) Naphthalene	10.381	128	22800	0.389	ng/ul	99
7) 2-Methylnaphthalene	12.003	142	14442	0.404	ng/ul	99
8) 1-Methylnaphthalene	12.229	142	14756	0.405	ng/ul	99
10) Acenaphthylene	13.919	152	20963	0.387	ng/ul	98
11) Acenaphthene	14.266	153	15868	0.387	ng/ul	99
12) Fluorene	15.261	166	17628	0.387	ng/ul	99
14) Pentachlorophenol	16.609	266	1552	0.303	ng/ul	98
15) Phenanthrene	16.993	178	26263	0.385	ng/ul	99
16) Anthracene	17.086	178	23341	0.384	ng/ul	99
19) Fluoranthene	19.011	202	27610	0.448	ng/ul	95
20) Pyrene	19.378	202	28111	0.423	ng/ul	96
21) Benzo(a)anthracene	21.131	228	19660	0.376	ng/ul	98
22) Chrysene	21.184	228	22210	0.375	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	18770	0.378	ng/ul	97
25) Benzo(k)fluoranthene	22.706	252	21846	0.380	ng/ul	96
26) Benzo(a)pyrene	23.212	252	17674	0.385	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.447	276	23838	0.370	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.467	278	18144	0.362	ng/ul#	95
29) Benzo(g,h,i)perylene	26.097	276	18805	0.362	ng/ul	97

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

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