

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032922\
 Data File : BM034399.D
 Acq On : 29 Mar 2022 10:16
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4062

Quant Time: Mar 30 02:28:19 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM032822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 28 18:27:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	2911	0.400	ng/ul	0.00
4) Naphthalene-d8	10.752	136	8220	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.573	164	4090	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.324	188	6512	0.400	ng/ul	0.00
17) Chrysene-d12	21.501	240	5131	0.400	ng/ul #	0.00
23) Perylene-d12	23.912	264	5782	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	1928	0.445	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152	4311	0.389	ng/ul	0.00
18) Fluoranthene-d10	19.348	212	6908	0.432	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.343	88	2377	0.497	ng/ul#	77
5) Naphthalene	10.796	128	9527	0.400	ng/ul#	90
7) 2-Methylnaphthalene	12.418	142	5727	0.393	ng/ul	98
8) 1-Methylnaphthalene	12.627	142	5665	0.390	ng/ul	100
10) Acenaphthylene	14.300	152	8061	0.410	ng/ul#	91
11) Acenaphthene	14.638	153	6154	0.411	ng/ul	96
12) Fluorene	15.628	166	6187	0.387	ng/ul	97
14) Pentachlorophenol	16.991	266	988	0.377	ng/ul	100
15) Phenanthrene	17.367	178	8393	0.403	ng/ul	97
16) Anthracene	17.464	178	7185	0.394	ng/ul	96
19) Fluoranthene	19.375	202	9600	0.415	ng/ul	97
20) Pyrene	19.738	202	10829	0.424	ng/ul#	94
21) Benzo(a)anthracene	21.486	228	5898	0.368	ng/ul	98
22) Chrysene	21.533	228	7335	0.400	ng/ul	98
24) Benzo(b)fluoranthene	23.173	252	7296	0.365	ng/ul#	73
25) Benzo(k)fluoranthene	23.222	252	7279	0.376	ng/ul#	81
26) Benzo(a)pyrene	23.810	252	8226	0.397	ng/ul#	66
27) Indeno(1,2,3-cd)pyrene	26.424	276	10853	0.402	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.451	278	8049	0.397	ng/ul#	73
29) Benzo(g,h,i)perylene	27.189	276	10722	0.411	ng/ul#	93

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

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