

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062322\
 Data File : BM035654.D
 Acq On : 24 Jun 2022 09:08
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4110

Quant Time: Jun 25 04:05:19 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.884	152	4124	0.400	ng/ul	0.03
4) Naphthalene-d8	10.644	136	18147	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.469	164	10491	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.229	188	21474	0.400	ng/ul	0.00
17) Chrysene-d12	21.409	240	16605	0.400	ng/ul	0.00
23) Perylene-d12	23.756	264	14429	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.268	96	2688	0.463	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.255	152	11064	0.410	ng/ul	0.01
18) Fluoranthene-d10	19.247	212	23233	0.417	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.302	88	2628	0.455	ng/ul#	59
5) Naphthalene	10.699	128	19421	0.316	ng/ul	99
7) 2-Methylnaphthalene	12.332	142	11523	0.328	ng/ul	93
8) 1-Methylnaphthalene	12.519	142	12529	0.389	ng/ul	100
10) Acenaphthylene	14.200	152	21073	0.357	ng/ul	99
11) Acenaphthene	14.529	153	15806	0.361	ng/ul	99
12) Fluorene	15.542	166	17772	0.359	ng/ul	100
14) Pentachlorophenol	16.933	266	1460	0.354	ng/ul	98
15) Phenanthrene	17.271	178	27231	0.346	ng/ul	98
16) Anthracene	17.381	178	23005	0.334	ng/ul	99
19) Fluoranthene	19.280	202	30373	0.358	ng/ul	98
20) Pyrene	19.642	202	33883	0.381	ng/ul	98
21) Benzo(a)anthracene	21.400	228	20124	0.320	ng/ul	99
22) Chrysene	21.447	228	25612	0.343	ng/ul	100
24) Benzo(b)fluoranthene	23.043	252	21377	0.338	ng/ul	99
25) Benzo(k)fluoranthene	23.092	252	21101	0.322	ng/ul	98
26) Benzo(a)pyrene	23.662	252	22098	0.337	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.208	276	25486	0.338	ng/ul	98
28) Dibenzo(a,h)anthracene	26.222	278	20591	0.334	ng/ul	97
29) Benzo(g,h,i)perylene	26.940	276	22881	0.339	ng/ul	97

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

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