

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
 Data File : BM035739.D
 Acq On : 29 Jun 2022 00:59
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4116

Quant Time: Jun 29 05:18:30 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 29 05:17:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	2487	0.400	ng/ul	0.00
4) Naphthalene-d8	10.622	136	10071	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.432	164	5274	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.196	188	9101	0.400	ng/ul	0.00
17) Chrysene-d12	21.373	240	10923	0.400	ng/ul	# 0.00
23) Perylene-d12	23.697	264	10887	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	1865	0.553	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.228	152	5587	0.354	ng/ul	0.00
18) Fluoranthene-d10	19.216	212	11222	0.305	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.251	88	1666	0.499	ng/ul#	86
5) Naphthalene	10.671	128	10636	0.345	ng/ul	98
7) 2-Methylnaphthalene	12.299	142	5842	0.303	ng/ul	99
8) 1-Methylnaphthalene	12.492	142	6363	0.349	ng/ul	99
10) Acenaphthylene	14.164	152	9954	0.360	ng/ul	98
11) Acenaphthene	14.497	153	7492	0.352	ng/ul	99
12) Fluorene	15.506	166	8078	0.329	ng/ul	99
14) Pentachlorophenol	16.896	266	849	0.591	ng/ul	98
15) Phenanthrene	17.238	178	11175	0.339	ng/ul	97
16) Anthracene	17.348	178	9433	0.344	ng/ul	97
19) Fluoranthene	19.248	202	14270	0.261	ng/ul	97
20) Pyrene	19.611	202	16190	0.264	ng/ul	97
21) Benzo(a)anthracene	21.362	228	12311	0.326	ng/ul	98
22) Chrysene	21.408	228	15609	0.321	ng/ul	99
24) Benzo(b)fluoranthene	22.989	252	15468	0.331	ng/ul	89
25) Benzo(k)fluoranthene	23.039	252	16557	0.347	ng/ul#	90
26) Benzo(a)pyrene	23.600	252	16870	0.354	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.107	276	21081	0.406	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.124	278	16983	0.413	ng/ul#	89
29) Benzo(g,h,i)perylene	26.845	276	19242	0.402	ng/ul	95

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

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