

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102422\
 Data File : BM037211.D
 Acq On : 24 Oct 2022 19:38
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4161

Quant Time: Oct 25 02:38:37 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 22 03:34:37 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	3965	0.400	ng/ul	0.00
4) Naphthalene-d8	10.699	136	12679	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.529	164	7497	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.271	188	16198	0.400	ng/ul	0.00
17) Chrysene-d12	21.447	240	14398	0.400	ng/ul	0.00
23) Perylene-d12	23.818	264	11966	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	1658	0.343	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.288	152	7163	0.397	ng/ul	0.00
18) Fluoranthene-d10	19.294	212	16779	0.385	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	1563	0.311	ng/ul#	80
5) Naphthalene	10.754	128	13808	0.368	ng/ul	100
7) 2-Methylnaphthalene	12.360	142	8896	0.396	ng/ul	98
8) 1-Methylnaphthalene	12.580	142	9176	0.400	ng/ul	99
10) Acenaphthylene	14.251	152	13344	0.414	ng/ul	99
11) Acenaphthene	14.593	153	10479	0.360	ng/ul	97
12) Fluorene	15.579	166	11832	0.348	ng/ul	100
14) Pentachlorophenol	16.925	266	1906	0.486	ng/ul	99
15) Phenanthrene	17.313	178	19950	0.365	ng/ul	99
16) Anthracene	17.402	178	17583	0.406	ng/ul	98
19) Fluoranthene	19.326	202	24025	0.384	ng/ul	96
20) Pyrene	19.689	202	24930	0.393	ng/ul	98
21) Benzo(a)anthracene	21.433	228	22481	0.402	ng/ul	99
22) Chrysene	21.485	228	22667	0.353	ng/ul	99
24) Benzo(b)fluoranthene	23.095	252	21658	0.342	ng/ul	98
25) Benzo(k)fluoranthene	23.142	252	20584	0.336	ng/ul	98
26) Benzo(a)pyrene	23.712	252	18027	0.354	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.272	276	21698	0.305	ng/ul	96
28) Dibenzo(a,h)anthracene	26.292	278	17100	0.305	ng/ul	97
29) Benzo(g,h,i)perylene	27.027	276	18109	0.290	ng/ul	98

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

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