

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061423\
 Data File : BM040238.D
 Acq On : 14 Jun 2023 15:08
 Operator : MA/JU
 Sample : SST0.885
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8018

Quant Time: Jun 15 00:21:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 15 00:19:21 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.974	152	12103	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136	42577	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.610	164	23125	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	44764	0.400	ng/ul	0.00
17) Chrysene-d12	21.530	240	35120	0.400	ng/ul	0.00
23) Perylene-d12	23.959	264	30957	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.350	96	11783	0.786	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.373	152	42067	0.779	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	70309	0.802	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.384	88	12094	0.770	ng/ul#	87
5) Naphthalene	10.833	128	79180	0.785	ng/ul	99
7) 2-Methylnaphthalene	12.445	142	49267	0.786	ng/ul	99
8) 1-Methylnaphthalene	12.665	142	51537	0.785	ng/ul	99
10) Acenaphthylene	14.332	152	61959	0.766	ng/ul	99
11) Acenaphthene	14.675	153	53700	0.774	ng/ul	99
12) Fluorene	15.655	166	59776	0.774	ng/ul	99
14) Pentachlorophenol	17.003	266	6854	0.751	ng/ul	100
15) Phenanthrene	17.396	178	91354	0.776	ng/ul	99
16) Anthracene	17.489	178	78165	0.784	ng/ul	99
19) Fluoranthene	19.408	202	95264	0.796	ng/ul	99
20) Pyrene	19.770	202	96696	0.778	ng/ul	98
21) Benzo(a)anthracene	21.515	228	78683	0.759	ng/ul	99
22) Chrysene	21.568	228	84451	0.771	ng/ul	99
24) Benzo(b)fluoranthene	23.213	252	90944	0.797	ng/ul	96
25) Benzo(k)fluoranthene	23.263	252	85618	0.798	ng/ul	94
26) Benzo(a)pyrene	23.848	252	73670	0.788	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.471	276	96141	0.811	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.494	278	77487	0.815	ng/ul	97
29) Benzo(g,h,i)perylene	27.249	276	79167	0.800	ng/ul	99

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

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