

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032039.D
 Acq On : 03 Sep 2021 13:56
 Operator : CG/JU
 Sample : SST0.424
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.4024

Quant Time: Sep 03 14:39:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM090321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 14:29:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.459	152	1163	0.400	ng/ul	0.00
4) Naphthalene-d8	10.221	136	3906	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.102	164	2382	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.865	188	4543	0.400	ng/ul	0.00
17) Chrysene-d12	21.074	240	3421	0.400	ng/ul	0.00
23) Perylene-d12	23.192	264	3321	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.071	96	546	0.423	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.822	152	2394	0.364	ng/ul	0.00
18) Fluoranthene-d10	18.904	212	4124	0.361	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.102	88	496	0.380	ng/ul	100
5) Naphthalene	10.271	128	4075	0.350	ng/ul	100
7) 2-Methylnaphthalene	11.899	142	2833	0.356	ng/ul	100
8) 1-Methylnaphthalene	12.115	142	2851	0.363	ng/ul	100
10) Acenaphthylene	13.824	152	3903	0.385	ng/ul	100
11) Acenaphthene	14.167	153	3125	0.366	ng/ul	100
12) Fluorene	15.171	166	3562	0.362	ng/ul	100
14) Pentachlorophenol	16.539	266	451	0.310	ng/ul	96
15) Phenanthrene	16.907	178	5503	0.381	ng/ul	100
16) Anthracene	17.008	178	4921	0.364	ng/ul	100
19) Fluoranthene	18.934	202	5957	0.408	ng/ul	100
20) Pyrene	19.300	202	6013	0.406	ng/ul	100
21) Benzo(a)anthracene	21.059	228	4855	0.351	ng/ul	100
22) Chrysene	21.110	228	5738	0.401	ng/ul	100
24) Benzo(b)fluoranthene	22.565	252	5878	0.393	ng/ul	100
25) Benzo(k)fluoranthene	22.609	252	6360	0.411	ng/ul	100
26) Benzo(a)pyrene	23.105	252	5339	0.407	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.271	276	7451	0.440	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.286	278	5968	0.437	ng/ul	100
29) Benzo(g,h,i)perylene	25.904	276	6613	0.458	ng/ul	100

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

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