

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083022\
 Data File : BM036454.D
 Acq On : 29 Aug 2022 15:13
 Operator : CG/JU
 Sample : SST0.463
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.4017

Quant Time: Aug 29 16:18:31 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM083022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 29 16:16:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.746	152	3999	0.400	ng/ul	0.00
4) Naphthalene-d8	10.541	136	14971	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.391	164	8410	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.133	188	18559	0.400	ng/ul	0.00
17) Chrysene-d12	21.321	240	16811	0.400	ng/ul	# 0.00
23) Perylene-d12	23.621	264	14159	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	2411	0.448	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.141	152	9464	0.395	ng/ul	0.00
18) Fluoranthene-d10	19.165	212	20497	0.987	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.181	88	2162	0.425	ng/ul	93
5) Naphthalene	10.590	128	16255	0.385	ng/ul	97
7) 2-Methylnaphthalene	12.212	142	10094	0.399	ng/ul	93
8) 1-Methylnaphthalene	12.432	142	10278	0.400	ng/ul	100
10) Acenaphthylene	14.113	152	12810	0.368	ng/ul	99
11) Acenaphthene	14.451	153	11368	0.404	ng/ul	99
12) Fluorene	15.441	166	12827	0.389	ng/ul	98
14) Pentachlorophenol	16.803	266	1322	0.360	ng/ul	99
15) Phenanthrene	17.175	178	21685	0.327	ng/ul	99
16) Anthracene	17.272	178	18373	0.349	ng/ul	99
19) Fluoranthene	19.197	202	23549	0.401	ng/ul	98
20) Pyrene	19.560	202	24357	0.399	ng/ul	99
21) Benzo(a)anthracene	21.306	228	20068	0.288	ng/ul	99
22) Chrysene	21.359	228	22662	0.237	ng/ul	99
24) Benzo(b)fluoranthene	22.919	252	22009	0.214	ng/ul	92
25) Benzo(k)fluoranthene	22.966	252	22021	0.212	ng/ul#	89
26) Benzo(a)pyrene	23.519	252	17815	0.273	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.993	276	21274	0.432	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.014	278	18265	0.437	ng/ul#	90
29) Benzo(g,h,i)perylene	26.718	276	20463	0.459	ng/ul	97

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

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