

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083022\
 Data File : BM036455.D
 Acq On : 29 Aug 2022 15:50
 Operator : CG/JU
 Sample : SST0.864
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8018

Quant Time: Aug 29 16:34:58 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	3239	0.400	ng/ul	0.00
4) Naphthalene-d8	10.541	136	12476	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.387	164	7095	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.133	188	15215	0.400	ng/ul	0.00
17) Chrysene-d12	21.318	240	13630	0.400	ng/ul	0.00
23) Perylene-d12	23.621	264	11344	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	3851	0.883	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.135	152	16142	0.809	ng/ul	0.00
18) Fluoranthene-d10	19.165	212	33415	1.985	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.181	88	3745	0.909	ng/ul#	90
5) Naphthalene	10.590	128	27097	0.771	ng/ul	96
7) 2-Methylnaphthalene	12.212	142	17132	0.812	ng/ul	95
8) 1-Methylnaphthalene	12.427	142	17646	0.825	ng/ul	100
10) Acenaphthylene	14.109	152	21891	0.745	ng/ul	98
11) Acenaphthene	14.451	153	19244	0.811	ng/ul	99
12) Fluorene	15.441	166	21459	0.772	ng/ul	99
14) Pentachlorophenol	16.799	266	2220	0.738	ng/ul	97
15) Phenanthrene	17.175	178	35655	0.657	ng/ul	99
16) Anthracene	17.268	178	30696	0.711	ng/ul	98
19) Fluoranthene	19.193	202	38687	0.812	ng/ul	99
20) Pyrene	19.555	202	39211	0.792	ng/ul	100
21) Benzo(a)anthracene	21.304	228	33486	0.593	ng/ul	98
22) Chrysene	21.353	228	36551	0.471	ng/ul	99
24) Benzo(b)fluoranthene	22.914	252	36662	0.444	ng/ul	88
25) Benzo(k)fluoranthene	22.963	252	37006	0.445	ng/ul#	85
26) Benzo(a)pyrene	23.513	252	30552	0.585	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	25.983	276	36566	0.926	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.004	278	31359	0.937	ng/ul#	87
29) Benzo(g,h,i)perylene	26.711	276	33921	0.950	ng/ul	97

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

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