

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101421\
 Data File : BM032507.D
 Acq On : 14 Oct 2021 15:59
 Operator : CG/JU
 Sample : SST0.126
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.1036

Quant Time: Oct 14 17:50:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 14 17:49:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.560	152	3918	0.400	ng/ul	0.00
4) Naphthalene-d8	10.348	136	11467	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.221	164	6054	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.957	188	11692	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.135	240	8686	0.400	ng/ul	0.00
23) Perylene-d12	23.276	264	8253	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.927	96	514	0.122	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.945	152	1541	0.101	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	2916	0.124	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.965	88	521	0.120	ng/ul#	58
5) Naphthalene	10.398	128	3657	0.108	ng/ul#	93
7) 2-Methylnaphthalene	12.017	142	2295	0.105	ng/ul	99
8) 1-Methylnaphthalene	12.242	142	2268	0.107	ng/ul	98
10) Acenaphthylene	13.937	152	2771	0.114	ng/ul	94
11) Acenaphthene	14.281	153	2468	0.111	ng/ul	99
12) Fluorene	15.267	166	2763	0.109	ng/ul#	94
14) Pentachlorophenol	16.627	266	254	0.086	ng/ul	97
15) Phenanthrene	16.998	178	4212	0.111	ng/ul	95
16) Anthracene	17.088	178	3588	0.114	ng/ul#	92
19) Fluoranthene	19.011	202	4775	0.132	ng/ul#	95
20) Pyrene	19.374	202	5038	0.139	ng/ul	95
21) Benzo(a)anthracene	21.121	228	3639	0.118	ng/ul	96
22) Chrysene	21.172	228	4109	0.114	ng/ul	98
24) Benzo(b)fluoranthene	22.641	252	3871	0.090	ng/ul	63
25) Benzo(k)fluoranthene	22.682	252	3852	0.093	ng/ul#	62
26) Benzo(a)pyrene	23.184	252	3234	0.101	ng/ul#	55
27) Indeno(1,2,3-cd)pyrene	25.384	276	4282	0.104	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.396	278	3425	0.102	ng/ul#	76
29) Benzo(g,h,i)perylene	26.038	276	3799	0.098	ng/ul#	89

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

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