

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072821\
 Data File : BM031205.D
 Acq On : 28 Jul 2021 11:03
 Operator : CG/JU
 Sample : SST0.417
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.4017

Quant Time: Jul 28 11:39:01 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 28 11:36:18 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.614	152	1051	0.400	ng/ul	0.00
4) Naphthalene-d8	10.379	136	3640	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.246	164	2284	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.992	188	4847	0.400	ng/ul	0.00
17) Chrysene-d12	21.181	240	4214	0.400	ng/ul	0.00
23) Perylene-d12	23.350	264	3960	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.203	96	429	0.389	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.975	152	2285	0.357	ng/ul	0.00
18) Fluoranthene-d10	19.024	212	5133	0.399	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.234	88	466	0.432	ng/ul	100
5) Naphthalene	10.428	128	4086	0.376	ng/ul	100
7) 2-Methylnaphthalene	12.047	142	2797	0.357	ng/ul	100
8) 1-Methylnaphthalene	12.268	142	2729	0.357	ng/ul	100
10) Acenaphthylene	13.968	152	3602	0.386	ng/ul	100
11) Acenaphthene	14.307	153	3038	0.368	ng/ul	100
12) Fluorene	15.301	166	3547	0.365	ng/ul	100
14) Pentachlorophenol	16.649	266	577	0.344	ng/ul	100
15) Phenanthrene	17.034	178	5751	0.374	ng/ul	100
16) Anthracene	17.124	178	5351	0.371	ng/ul	100
19) Fluoranthene	19.054	202	6563	0.387	ng/ul	100
20) Pyrene	19.416	202	6686	0.386	ng/ul	100
21) Benzo(a)anthracene	21.166	228	6274	0.373	ng/ul	100
22) Chrysene	21.217	228	6587	0.362	ng/ul	100
24) Benzo(b)fluoranthene	22.706	252	6627	0.356	ng/ul	100
25) Benzo(k)fluoranthene	22.747	252	6838	0.362	ng/ul	100
26) Benzo(a)pyrene	23.256	252	5736	0.356	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.497	276	7644	0.363	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.514	278	6161	0.364	ng/ul	100
29) Benzo(g,h,i)perylene	26.149	276	6564	0.367	ng/ul	100

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

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