

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071223\
 Data File : BN026477.D
 Acq On : 12 Jul 2023 16:44
 Operator : MA/JU
 Sample : SST0.415
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4232

Quant Time: Jul 13 04:32:57 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 04:32:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.014	152	2230	0.400	ng/ul	0.00
4) Naphthalene-d8	10.809	136	7666	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.626	164	4231	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.372	188	6470	0.400	ng/ul	0.00
17) Chrysene-d12	21.564	240	5227	0.400	ng/ul	0.00
23) Perylene-d12	24.002	264	6251	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	1284	0.437	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.398	152	4089	0.392	ng/ul	0.00
18) Fluoranthene-d10	19.401	212	6783	0.374	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.381	88	1231	0.443	ng/ul	100
5) Naphthalene	10.858	128	8148	0.387	ng/ul	100
7) 2-Methylnaphthalene	12.469	142	4637	0.388	ng/ul	100
8) 1-Methylnaphthalene	12.684	142	5150	0.403	ng/ul	100
10) Acenaphthylene	14.353	152	7516	0.363	ng/ul	100
11) Acenaphthene	14.686	153	5866	0.372	ng/ul	100
12) Fluorene	15.680	166	5837	0.352	ng/ul	100
14) Pentachlorophenol	17.043	266	635	0.389	ng/ul	100
15) Phenanthrene	17.414	178	7666	0.372	ng/ul	100
16) Anthracene	17.516	178	6564	0.367	ng/ul	100
19) Fluoranthene	19.428	202	8904	0.363	ng/ul	100
20) Pyrene	19.796	202	9599	0.361	ng/ul	100
21) Benzo(a)anthracene	21.546	228	7135	0.371	ng/ul	100
22) Chrysene	21.602	228	8120	0.371	ng/ul	100
24) Benzo(b)fluoranthene	23.259	252	8849	0.363	ng/ul	100
25) Benzo(k)fluoranthene	23.309	252	9421	0.369	ng/ul	100
26) Benzo(a)pyrene	23.897	252	8020	0.370	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.557	276	11154	0.393	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.578	278	8460	0.388	ng/ul	100
29) Benzo(g,h,i)perylene	27.346	276	9972	0.398	ng/ul	100

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

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