

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111321\
 Data File : BN017398.D
 Acq On : 13 Nov 2021 18:59
 Operator : CG/JU
 Sample : SST0.839
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8239

Quant Time: Nov 15 02:23:11 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 02:20:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.877	152	6431	0.400	ng/ul	0.00
4) Naphthalene-d8	10.674	136	23899	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.509	164	15271	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.253	188	32702	0.400	ng/ul	0.00
17) Chrysene-d12	21.437	240	26280	0.400	ng/ul	0.00
23) Perylene-d12	23.819	264	20342	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.333	96	5563	0.778	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.264	152	28473	0.816	ng/ul	0.00
18) Fluoranthene-d10	19.283	212	69653	1.011	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.366	88	6846	0.891	ng/ul#	78
5) Naphthalene	10.724	128	51563	0.779	ng/ul	97
7) 2-Methylnaphthalene	12.335	142	36634	0.800	ng/ul	99
8) 1-Methylnaphthalene	12.555	142	36198	0.772	ng/ul	100
10) Acenaphthylene	14.226	152	56763	0.940	ng/ul	99
11) Acenaphthene	14.573	153	40425	0.763	ng/ul	99
12) Fluorene	15.559	166	48001	0.775	ng/ul	98
14) Pentachlorophenol	16.903	266	5599	0.759	ng/ul	99
15) Phenanthrene	17.295	178	75863	0.731	ng/ul	99
16) Anthracene	17.384	178	76659	0.891	ng/ul	99
19) Fluoranthene	19.311	202	89596	0.953	ng/ul	100
20) Pyrene	19.673	202	91699	0.951	ng/ul	97
21) Benzo(a)anthracene	21.422	228	75505	0.877	ng/ul	100
22) Chrysene	21.475	228	75048	0.769	ng/ul	100
24) Benzo(b)fluoranthene	23.097	252	64026	0.844	ng/ul	90
25) Benzo(k)fluoranthene	23.144	252	65891	0.766	ng/ul#	90
26) Benzo(a)pyrene	23.714	252	53776	0.791	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.284	276	46100	0.478	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.307	278	34643	0.454	ng/ul#	89
29) Benzo(g,h,i)perylene	27.035	276	41432	0.507	ng/ul	94

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

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