

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090822\
 Data File : BN021640.D
 Acq On : 08 Sep 2022 11:38
 Operator : CG/JU
 Sample : SST0.442
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4246

Quant Time: Sep 08 13:40:58 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 08 13:38:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.067	152	4689	0.400	ng/u1	0.00
4) Naphthalene-d8	10.894	136	15043	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.721	164	8584	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.468	188	17675	0.400	ng/u1	0.00
17) Chrysene-d12	21.662	240	12048	0.400	ng/u1	0.00
23) Perylene-d12	24.176	264	8079	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	2359	0.390	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.483	152	9641	0.424	ng/u1	0.00
18) Fluoranthene-d10	19.497	212	18041	0.486	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	2469	0.390	ng/u1	93
5) Naphthalene	10.944	128	17220	0.362	ng/u1	98
7) 2-Methylnaphthalene	12.555	142	11120	0.373	ng/u1	98
8) 1-Methylnaphthalene	12.775	142	11483	0.411	ng/u1	99
10) Acenaphthylene	14.444	152	14472	0.327	ng/u1	98
11) Acenaphthene	14.781	153	12437	0.367	ng/u1	98
12) Fluorene	15.767	166	13939	0.368	ng/u1	100
14) Pentachlorophenol	17.118	266	1272	0.652	ng/u1	99
15) Phenanthrene	17.510	178	22447	0.343	ng/u1	99
16) Anthracene	17.603	178	18363	0.350	ng/u1	99
19) Fluoranthene	19.524	202	22639	0.396	ng/u1	96
20) Pyrene	19.892	202	22267	0.390	ng/u1	98
21) Benzo(a)anthracene	21.644	228	15796	0.378	ng/u1	100
22) Chrysene	21.697	228	17636	0.328	ng/u1	99
24) Benzo(b)fluoranthene	23.401	252	15605	0.450	ng/u1	98
25) Benzo(k)fluoranthene	23.451	252	14963	0.400	ng/u1	98
26) Benzo(a)pyrene	24.065	252	12074	0.369	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.830	276	13555	0.353	ng/u1#	95
28) Dibenzo(a,h)anthracene	26.851	278	11401	0.361	ng/u1	96
29) Benzo(g,h,i)perylene	27.639	276	11944	0.313	ng/u1	97

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

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