

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102022\
 Data File : BN022192.D
 Acq On : 19 Oct 2022 11:59
 Operator : CG/JU
 Sample : SST0.448
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 FID_A
 ClientSampleId :
 SST0.4204

Quant Time: Oct 19 15:03:58 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 00:20:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	0.000	152	0	0.00	ng/ul	-8.05
4) Naphthalene-d8	0.000	136	0	0.00	ng/ul	-10.88
9) Acenaphthene-d10	14.627	164	5296	0.40	ng/ul	-0.08
13) Phenanthrene-d10	17.382	188	15404	0.40	ng/ul	#-0.08
17) Chrysene-d12	21.582	240	12145	0.40	ng/ul	#-0.06
23) Perylene-d12	24.047	264	8849	0.40	ng/ul	#-0.11
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.421	96	20	0.00	ng/ul	0.06
6) 2-Methylnaphthalene-d10	0.000	152	0	0.00	ng/ul	
18) Fluoranthene-d10	19.415	212	16884	0.37	ng/ul	-0.07
Target Compounds						
						Qvalue
10) Acenaphthylene	14.349	152	12899	0.58	ng/ul	98
11) Acenaphthene	14.687	153	10948	0.58	ng/ul	100
12) Fluorene	15.677	166	12711	0.60	ng/ul	99
14) Pentachlorophenol	17.031	266	1319	0.44	ng/ul	99
15) Phenanthrene	17.424	178	21257	0.44	ng/ul	99
16) Anthracene	17.517	178	17620	0.44	ng/ul	98
19) Fluoranthene	19.443	202	22923	0.40	ng/ul	96
20) Pyrene	19.810	202	22747	0.41	ng/ul	94
21) Benzo(a)anthracene	21.568	228	18435	0.45	ng/ul	99
22) Chrysene	21.620	228	19825	0.45	ng/ul	99
24) Benzo(b)fluoranthene	23.289	252	18761	0.44	ng/ul	97
25) Benzo(k)fluoranthene	23.339	252	18557	0.45	ng/ul	96
26) Benzo(a)pyrene	23.936	252	14976	0.46	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.636	276	18407	0.50	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.656	278	14790	0.48	ng/ul	94
29) Benzo(g,h,i)perylene	27.430	276	15179	0.45	ng/ul	97

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

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