

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101922\
 Data File : BM037149.D
 Acq On : 19 Oct 2022 12:58
 Operator : CG/JU
 Sample : SST0.876
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8032

Quant Time: Oct 20 02:00:39 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 20 01:35:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.909	152	3555	0.400	ng/u1	0.00
4) Naphthalene-d8	10.710	136	10301	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.538	164	5534	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.280	188	12166	0.400	ng/u1	0.00
17) Chrysene-d12	21.456	240	10766	0.400	ng/u1	0.00
23) Perylene-d12	23.835	264	9312	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	3531	0.704	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.294	152	12475	0.802	ng/u1	0.00
18) Fluoranthene-d10	19.303	212	27027	0.839	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	3649	0.745	ng/u1	93
5) Naphthalene	10.760	128	25602	0.869	ng/u1	99
7) 2-Methylnaphthalene	12.371	142	15843	0.879	ng/u1	98
8) 1-Methylnaphthalene	12.585	142	15960	0.872	ng/u1	100
10) Acenaphthylene	14.261	152	19909	0.866	ng/u1	98
11) Acenaphthene	14.598	153	18058	0.905	ng/u1	100
12) Fluorene	15.584	166	21249	0.939	ng/u1	99
14) Pentachlorophenol	16.933	266	2385	0.810	ng/u1	98
15) Phenanthrene	17.322	178	34531	0.889	ng/u1	99
16) Anthracene	17.415	178	28032	0.864	ng/u1	99
19) Fluoranthene	19.331	202	39777	0.929	ng/u1	100
20) Pyrene	19.698	202	39555	0.886	ng/u1	99
21) Benzo(a)anthracene	21.441	228	35169	0.904	ng/u1	100
22) Chrysene	21.494	228	39748	0.952	ng/u1	99
24) Benzo(b)fluoranthene	23.110	252	41338	0.994	ng/u1	96
25) Benzo(k)fluoranthene	23.157	252	40610	1.002	ng/u1	94
26) Benzo(a)pyrene	23.730	252	33190	0.983	ng/u1#	91
27) Indeno(1,2,3-cd)pyrene	26.299	276	47536	0.994	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.319	278	37937	0.982	ng/u1	97
29) Benzo(g,h,i)perylene	27.057	276	41253	0.979	ng/u1	98

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

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