

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042222\
 Data File : BM034767.D
 Acq On : 22 Apr 2022 14:26
 Operator : CG/JU
 Sample : SST0.431
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.4031

Quant Time: Apr 22 16:10:14 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM042222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 22 16:08:48 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	4825	0.400	ng/ul	0.00
4) Naphthalene-d8	10.747	136	12345	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.574	164	6672	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	12993	0.400	ng/ul #	0.00
17) Chrysene-d12	21.486	240	8660	0.400	ng/ul #	0.00
23) Perylene-d12	23.883	264	7436	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.386	96	2387	0.353	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.336	152	7416	0.402	ng/ul	0.00
18) Fluoranthene-d10	19.339	212	12220	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.420	88	2204	0.407	ng/ul#	78
5) Naphthalene	10.797	128	14979	0.342	ng/ul#	88
7) 2-Methylnaphthalene	12.408	142	9990	0.364	ng/ul	100
8) 1-Methylnaphthalene	12.628	142	9803	0.437	ng/ul	99
10) Acenaphthylene	14.296	152	12173	0.301	ng/ul#	89
11) Acenaphthene	14.639	153	10336	0.340	ng/ul	96
12) Fluorene	15.619	166	11684	0.339	ng/ul#	98
14) Pentachlorophenol	16.962	266	1501	0.355	ng/ul	97
15) Phenanthrene	17.354	178	17977	0.337	ng/ul	94
16) Anthracene	17.447	178	15939	0.333	ng/ul#	91
19) Fluoranthene	19.366	202	18344	0.351	ng/ul	96
20) Pyrene	19.729	202	18090	0.349	ng/ul	99
21) Benzo(a)anthracene	21.472	228	14061	0.333	ng/ul	97
22) Chrysene	21.524	228	14959	0.327	ng/ul	97
24) Benzo(b)fluoranthene	23.152	252	14811	0.343	ng/ul	93
25) Benzo(k)fluoranthene	23.202	252	14076	0.342	ng/ul	93
26) Benzo(a)pyrene	23.775	252	11781	0.304	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.361	276	16054	0.326	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.384	278	12828	0.315	ng/ul#	89
29) Benzo(g,h,i)perylene	27.119	276	14193	0.315	ng/ul	95

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

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