

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082622\
 Data File : BM036422.D
 Acq On : 26 Aug 2022 12:39
 Operator : CG/JU
 Sample : SST0.859
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8011

Quant Time: Aug 26 13:29:34 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM082622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 13:25:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.795	152	3973	0.400	ng/ul	0.00
4) Naphthalene-d8	10.589	136	16052	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.432	164	9139	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.179	188	21394	0.400	ng/ul	0.00
17) Chrysene-d12	21.362	240	45580	0.400	ng/ul #	0.00
23) Perylene-d12	23.691	264	22863	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	4423	0.820	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.184	152	21119	0.840	ng/ul	0.00
18) Fluoranthene-d10	19.206	212	49317	0.321	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	4231	0.793	ng/ul#	87
5) Naphthalene	10.639	128	34299	0.699	ng/ul	97
7) 2-Methylnaphthalene	12.261	142	22064	0.718	ng/ul	93
8) 1-Methylnaphthalene	12.475	142	22517	0.774	ng/ul	99
10) Acenaphthylene	14.155	152	30624	0.639	ng/ul	97
11) Acenaphthene	14.493	153	24844	0.673	ng/ul	100
12) Fluorene	15.487	166	28676	0.674	ng/ul	99
14) Pentachlorophenol	16.846	266	3323	0.984	ng/ul	98
15) Phenanthrene	17.221	178	48781	0.629	ng/ul	99
16) Anthracene	17.314	178	44058	0.684	ng/ul	98
19) Fluoranthene	19.234	202	70353	0.308	ng/ul	97
20) Pyrene	19.597	202	73218	0.286	ng/ul	97
21) Benzo(a)anthracene	21.344	228	78907	0.500	ng/ul	98
22) Chrysene	21.397	228	97682	0.482	ng/ul	98
24) Benzo(b)fluoranthene	22.975	252	79351	0.810	ng/ul	85
25) Benzo(k)fluoranthene	23.022	252	82847	0.827	ng/ul#	82
26) Benzo(a)pyrene	23.586	252	58379	0.583	ng/ul#	77
27) Indeno(1,2,3-cd)pyrene	26.094	276	59829	0.548	ng/ul	93
28) Dibenzo(a,h)anthracene	26.117	278	51486	0.596	ng/ul#	85
29) Benzo(g,h,i)perylene	26.828	276	55470	0.551	ng/ul	92

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

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