

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033177.D
 Acq On : 19 Nov 2021 15:53
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV014

Quant Time: Nov 19 16:29:20 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 15:41:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	2416	0.400	ng/ul	0.00
4) Naphthalene-d8	10.760	136	6637	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.580	164	4014	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	8888	0.400	ng/ul	0.00
17) Chrysene-d12	21.475	240	7616	0.400	ng/ul	0.00
23) Perylene-d12	23.817	264	5852	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.417	96	901	0.397	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	3766	0.409	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	8863	0.402	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.451	88	1018	0.398	ng/ul#	92
5) Naphthalene	10.810	128	8101	0.394	ng/ul	99
7) 2-Methylnaphthalene	12.415	142	5408	0.406	ng/ul	98
8) 1-Methylnaphthalene	12.631	142	5436	0.413	ng/ul	99
10) Acenaphthylene	14.303	152	7337	0.391	ng/ul#	100
11) Acenaphthene	14.644	153	6116	0.397	ng/ul	98
12) Fluorene	15.626	166	7141	0.411	ng/ul	97
14) Pentachlorophenol	16.969	266	1039	0.350	ng/ul	97
15) Phenanthrene	17.358	178	11692	0.363	ng/ul	99
16) Anthracene	17.448	178	10164	0.390	ng/ul	99
19) Fluoranthene	19.365	202	13506	0.368	ng/ul	97
20) Pyrene	19.723	202	13419	0.369	ng/ul	97
21) Benzo(a)anthracene	21.458	228	10650	0.392	ng/ul	99
22) Chrysene	21.509	228	11675	0.390	ng/ul	99
24) Benzo(b)fluoranthene	23.103	252	11040	0.417	ng/ul	98
25) Benzo(k)fluoranthene	23.153	252	10828	0.414	ng/ul	98
26) Benzo(a)pyrene	23.711	252	8686	0.396	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.213	276	10792	0.390	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.228	278	8492	0.384	ng/ul	95
29) Benzo(g,h,i)perylene	26.955	276	9528	0.383	ng/ul	94

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

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