

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM092920\
 Data File : BM027569.D
 Acq On : 29 Sep 2020 15:10
 Operator : JU/CG
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV001

Quant Time: Sep 29 15:41:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM092920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 29 15:04:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	459	0.40	ng/ul	0.00
2) Naphthalene-d8	10.59	136	1208	0.40	ng/ul	0.00
7) Acenaphthene-d10	14.43	164	915	0.40	ng/ul	0.00
11) Phenanthrene-d10	17.17	188	2400	0.40	ng/ul	0.00
17) Chrysene-d12	21.36	240	2633	0.40	ng/ul	0.00
21) Perylene-d12	23.62	264	2732	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.18	152	916	0.38	ng/ul	0.00
15) Fluoranthene-d10	19.20	212	3064	0.37	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.64	128	1354	0.404	ng/ul	98
5) 2-Methylnaphthalene	12.25	142	1119	0.432	ng/ul	99
6) 1-Methylnaphthalene	12.47	142	1011	0.342	ng/ul	99
8) Acenaphthylene	14.15	152	1689	0.410	ng/ul	99
9) Acenaphthene	14.50	153	1237	0.402	ng/ul	99
10) Fluorene	15.48	166	1641	0.405	ng/ul	97
12) Pentachlorophenol	16.83	266	376	0.401	ng/ul	100
13) Phenanthrene	17.22	178	2852	0.399	ng/ul	99
14) Anthracene	17.31	178	2745	0.400	ng/ul	99
16) Fluoranthene	19.23	202	4088	0.406	ng/ul	99
18) Pyrene	19.59	202	4025	0.403	ng/ul	99
19) Benzo(a)anthracene	21.34	228	4246	0.410	ng/ul	99
20) Chrysene	21.39	228	4470	0.421	ng/ul	100
22) Benzo(b)fluoranthene	22.94	252	4684	0.416	ng/ul	98
23) Benzo(k)fluoranthene	22.99	252	4617	0.416	ng/ul	97
24) Benzo(a)pyrene	23.52	252	4176	0.415	ng/ul	99
25) Indeno(1,2,3-cd)pyrene	25.90	276	5610	0.420	ng/ul#	99
26) Dibenzo(a,h)anthracene	25.92	278	4666	0.420	ng/ul	98
27) Benzo(a,h,i)perylene	26.60	276	4684	0.408	ng/ul	98

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

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