

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090718\
 Data File : BF108864.D
 Acq On : 7 Sep 2018 16:52
 Operator : JU/SJ
 Sample : J4779-01 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-090618-A

Manual Integrations
 APPROVED

Sohil
 9/10/2018 10:39:31 AM

Quant Time: Sep 07 18:59:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	245760	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	945466	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	420853	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	693044	20.00	ng	0.00
75) Chrysene-d12	13.87	240	517784	20.00	ng	0.00
86) Perylene-d12	15.28	264	477907	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	296241	19.97	ng	0.00
7) Phenol-d6	6.36	99	390345	20.45	ng	0.00
23) Nitrobenzene-d5	7.29	82	227834	15.08	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	69286	17.25	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	436225	17.68	ng	0.00
78) Terphenyl-d14	12.82	244	315358	14.20	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	98513	3.347	ng	Qvalue # 99
57) Fluorene	10.31	166	87234	3.858	ng	99
70) Phenanthrene	11.27	178	601724	18.181	ng	99
71) Anthracene	11.32	178	195340	6.165	ng	98
72) Carbazole	11.47	167	89725	2.689	ng	99
74) Fluoranthene	12.45	202	760207	23.200	ng	96
77) Pyrene	12.68	202	694290	17.426	ng	99
80) Benzo(a)anthracene	13.86	228	362766	10.931	ng	98
82) Chrysene	13.90	228	303353	9.442	ng	98
87) Benzo(b)fluoranthene	14.88	252	386830m	14.886	ng	
88) Benzo(k)fluoranthene	14.90	252	123746m	5.171	ng	
89) Benzo(a)pyrene	15.22	252	272336	11.615	ng	98
91) Benzo(a,h,i)perylene	17.04	276	123822	5.985	ng	94

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

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