

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071118\
 Data File : BF107301.D
 Acq On : 11 Jul 2018 15:45
 Operator : JU/SJ
 Sample : J3933-03 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-3-6

Manual Integrations
 APPROVED

Sohil
 7/12/2018 2:19:21 PM

Quant Time: Jul 11 16:59:49 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 11 16:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	58122	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	209363	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	87100	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	151804	20.00	ng	0.00
75) Chrysene-d12	14.14	240	179618	20.00	ng	0.00
86) Perylene-d12	15.68	264	172211	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	89411	26.05	ng	0.00
7) Phenol-d6	6.58	99	110516	25.78	ng	0.00
23) Nitrobenzene-d5	7.50	82	64907	17.23	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	22041	21.83	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	109999	8.02	ng	0.00
78) Terphenyl-d14	13.06	244	112369	15.15	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	19538	2.991	ng	99
70) Phenanthrene	11.50	178	295361	40.340	ng	100
71) Anthracene	11.55	178	81103	10.852	ng	97
72) Carbazole	11.71	167	22973	3.283	ng	97
74) Fluoranthene	12.71	202	701742	86.956	ng	95
77) Pyrene	12.94	202	748314	63.178	ng	100
80) Benzo(a)anthracene	14.13	228	526618	48.105	ng	99
82) Chrysene	14.17	228	448935	44.575	ng	98
85) Indeno(1,2,3-cd)pyrene	17.28	276	182492m	21.352	ng	
87) Benzo(b)fluoranthene	15.23	252	603153m	56.382	ng	
88) Benzo(k)fluoranthene	15.25	252	154352m	15.151	ng	
89) Benzo(a)pyrene	15.62	252	409113m	43.019	ng	
90) Dibenzo(a,h)anthracene	17.27	278	48319m	5.995	ng	
91) Benzo(g,h,i)perylene	17.76	276	166894	21.186	ng	89

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

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