

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112862.D
 Acq On : 5 Mar 2019 19:50
 Operator : JU/SJ
 Sample : K1705-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EXT-028-SW009-022719

Manual Integrations
 APPROVED

Sohil
 3/6/2019 12:24:29 PM

Quant Time: Mar 06 03:53:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	174881	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	674296	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	317280	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	596150	20.00	ng	0.00
76) Chrysene-d12	13.87	240	460944	20.00	ng	0.00
87) Perylene-d12	15.28	264	484311	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1244381	110.69	ng	0.00
7) Phenol-d6	6.38	99	1609376	113.96	ng	0.01
23) Nitrobenzene-d5	7.30	82	1000375	88.77	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	448953	133.29	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1478750	74.82	ng	0.00
79) Terphenyl-d14	12.83	244	1250957	52.61	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	147164	6.290	ng	99
71) Phenanthrene	11.27	178	238305	8.034	ng	98
75) Fluoranthene	12.45	202	281061	8.844	ng	96
78) Pyrene	12.68	202	225827	5.811	ng	99
81) Benzo(a)anthracene	13.86	228	110669	3.464	ng	96
83) Chrysene	13.90	228	95301	3.045	ng	96
88) Benzo(b)fluoranthene	14.88	252	122376m	3.906	ng	
90) Benzo(a)pyrene	15.22	252	84789	3.060	ng	98

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

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