

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081618\
 Data File : BF108199.D
 Acq On : 16 Aug 2018 17:09
 Operator : JU/SJ
 Sample : J4499-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LSB-04(0-1)

Quant Time: Aug 17 14:40:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	181691	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	662384	20.00	ng	0.00
38) Acenaphthene-d10	10.06	164	296410	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	460214	20.00	ng	0.00
75) Chrysene-d12	14.21	240	431895	20.00	ng	0.00
86) Perylene-d12	15.77	264	404371	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	927822	92.45	ng	0.01
7) Phenol-d6	6.64	99	1229020	92.16	ng	0.00
23) Nitrobenzene-d5	7.58	82	831974	70.09	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	267280	90.40	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	1457371	78.94	ng	0.00
78) Terphenyl-d14	13.14	244	1133276	66.72	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.76	163	129378	6.266	ng	Qvalue # 99
70) Phenanthrene	11.58	178	54630	2.517	ng	99
74) Fluoranthene	12.78	202	148487	6.268	ng	93
77) Pyrene	13.01	202	128863	4.713	ng	97
80) Benzo(a)anthracene	14.20	228	72198	2.913	ng	96
82) Chrysene	14.24	228	79027	3.478	ng	97
87) Benzo(b)fluoranthene	15.30	252	112927m	4.674	ng	
89) Benzo(a)pyrene	15.70	252	69058	3.192	ng	95

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

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