

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080219\
 Data File : BF115965.D
 Acq On : 2 Aug 2019 23:38
 Operator : HP/JU
 Sample : K4048-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWL-C4-GW

Quant Time: Aug 05 16:38:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:31:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	105370	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	392631	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	226854	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	397201	20.00	ng	0.00
76) Chrysene-d12	14.02	240	279588	20.00	ng	0.00
87) Perylene-d12	15.50	264	319379	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	397133	63.09	ng	0.02
7) Phenol-d6	6.48	99	328980	41.43	ng	0.00
23) Nitrobenzene-d5	7.42	82	630801	92.74	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	294454	133.88	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1124415	92.84	ng	0.00
79) Terphenyl-d14	12.97	244	1349223	101.69	ng	0.00
Target Compounds						
10) Phenol	6.49	94	106089	11.344	ng	99
32) Benzoic acid	7.90	122	37710	10.269	ng	93
50) Dimethylphthalate	9.60	163	119210	7.718	ng	98
52) Acenaphthene	9.93	154	65976	5.530	ng	93
58) Fluorene	10.44	166	105605	7.911	ng	98
71) Phenanthrene	11.40	178	102643	5.055	ng	98

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

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