

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081019\
 Data File : BF116158.D
 Acq On : 10 Aug 2019 19:28
 Operator : HP/JU
 Sample : K4257-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAND-A

Quant Time: Aug 12 03:07:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	158422	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	589327	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	270023	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	389672	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	285240	20.00	ng	-0.01
87) Perylene-d12	15.47	264	272988	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1015554	107.31	ng	0.00
7) Phenol-d6	6.48	99	1308897	109.63	ng	0.00
23) Nitrobenzene-d5	7.40	82	860732	84.31	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	243460	93.00	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1403361	97.35	ng	-0.01
79) Terphenyl-d14	12.94	244	1124736	83.09	ng	-0.01
Target Compounds						
10) Phenol	6.49	94	33403	2.376	ng	91
50) Dimethylphthalate	9.59	163	205199	11.161	ng	99
71) Phenanthrene	11.38	178	40067	2.011	ng	97
75) Fluoranthene	12.57	202	54147	2.596	ng	99
78) Pyrene	12.80	202	47828	2.224	ng	97

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

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