

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082319\
 Data File : BF116517.D
 Acq On : 24 Aug 2019 12:19
 Operator : HP/JU
 Sample : K4490-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAND-STONE

Manual Integrations
 APPROVED

Jagrut
 8/28/2019 8:09:05 AM

Quant Time: Aug 26 03:54:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 23 13:38:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	128200	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	465088	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	223616	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	359297	20.00	ng	0.00
76) Chrysene-d12	14.03	240	261626	20.00	ng	0.00
87) Perylene-d12	15.49	264	240736	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	678011	77.28	ng	0.00
7) Phenol-d6	6.51	99	912700	84.84	ng	0.00
23) Nitrobenzene-d5	7.44	82	541620	63.71	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	174326	90.79	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	1045391	75.29	ng	0.00
79) Terphenyl-d14	12.98	244	1023026	73.04	ng	0.00
Target Compounds						
10) Phenol	6.52	94	37538	3.213	ng	85
50) Dimethylphthalate	9.62	163	149417	9.577	ng	99

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

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