

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
 Data File : BF111970.D
 Acq On : 10 Jan 2019 00:42
 Operator : JU/SJ
 Sample : K1075-15RE 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BASIN-8RE

Manual Integrations
 APPROVED

Sohil
 1/10/2019 11:46:10 AM

Quant Time: Jan 10 03:27:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	102815	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	427969	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	207956	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	340624	20.00	ng	0.00
76) Chrysene-d12	14.06	240	130397	20.00	ng	0.00
87) Perylene-d12	15.55	264	168981	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	148947	24.78	ng	0.00
7) Phenol-d6	6.52	99	165052	21.31	ng	-0.01
23) Nitrobenzene-d5	7.44	82	95520	11.93	ng	-0.02
42) 2,4,6-Tribromophenol	10.72	330	37232	13.76	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	209192	14.67	ng	-0.01
79) Terphenyl-d14	12.99	244	91641	13.34	ng	0.00
Target Compounds						
71) Phenanthrene	11.44	178	53283	2.926	ng	97
75) Fluoranthene	12.63	202	40963	2.191	ng	98
78) Pyrene	12.86	202	34084	3.799	ng	98
80) Butylbenzylphthalate	13.47	149	30259	7.977	ng	96
81) Benzo(a)anthracene	14.05	228	15034	2.003	ng	93
88) Benzo(b)fluoranthene	15.11	252	24646m	2.438	ng	

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

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