

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051319\
 Data File : BF114254.D
 Acq On : 14 May 2019 10:43
 Operator : JU/SJ
 Sample : K2765-21DL 4X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: May 14 12:26:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	257731	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	1050816	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	500403	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	928898	20.00	ng	0.00
76) Chrysene-d12	14.02	240	623859	20.00	ng	0.00
87) Perylene-d12	15.50	264	614752	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	193856	12.10	ng	0.00
7) Phenol-d6	6.50	99	273102	14.42	ng	0.00
23) Nitrobenzene-d5	7.41	82	159785	8.53	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	52133	10.08	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	239176	7.23	ng	0.00
79) Terphenyl-d14	12.95	244	197711	6.53	ng	0.00
Target Compounds						
49) Acenaphthylene	9.75	152	168647	3.408	ng	96
71) Phenanthrene	11.40	178	603011	13.343	ng	98
72) Anthracene	11.45	178	96981	2.137	ng	98
75) Fluoranthene	12.59	202	804307	16.527	ng	98
78) Pyrene	12.82	202	1479649	29.483	ng	99
81) Benzo(a)anthracene	14.01	228	487953m	12.093	ng	
83) Chrysene	14.04	228	591853	14.963	ng	97
86) Indeno(1,2,3-cd)pyrene	16.98	276	184141	5.267	ng	96
88) Benzo(b)fluoranthene	15.07	252	513485m	13.038	ng	
89) Benzo(k)fluoranthene	15.09	252	172481m	4.767	ng	
90) Benzo(a)pyrene	15.44	252	346662	10.001	ng	98
92) Benzo(a,h,i)perylene	17.43	276	204143	7.073	ng	97

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

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