

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030419\
 Data File : BF112823.D
 Acq On : 4 Mar 2019 14:45
 Operator : JU/SJ
 Sample : K1707-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S12

Manual Integrations
 APPROVED

Sohil
 3/5/2019 11:00:43 AM

Quant Time: Mar 05 01:17:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	160205	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	613328	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	269431	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	465139	20.00	ng	0.00
76) Chrysene-d12	13.88	240	401901	20.00	ng	0.00
87) Perylene-d12	15.29	264	401016	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	992549	96.37	ng	0.01
7) Phenol-d6	6.38	99	1307882	101.09	ng	0.01
23) Nitrobenzene-d5	7.30	82	843796	82.32	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	310331	108.49	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1263322	75.41	ng	0.00
79) Terphenyl-d14	12.83	244	963948	46.49	ng	0.00
Target Compounds						
10) Phenol	6.39	94	32028	2.122	ng	91
50) Dimethylphthalate	9.49	163	95418	4.803	ng	98
71) Phenanthrene	11.27	178	101885	4.403	ng	98
75) Fluoranthene	12.46	202	162234	6.543	ng	96
78) Pyrene	12.69	202	166302	4.908	ng	96
81) Benzo(a)anthracene	13.87	228	98437m	3.534	ng	
83) Chrysene	13.90	228	113116	4.146	ng	96
86) Indeno(1,2,3-cd)pyrene	16.65	276	48793	2.098	ng	95
88) Benzo(b)fluoranthene	14.89	252	150324m	5.795	ng	
90) Benzo(a)pyrene	15.23	252	93122	4.059	ng	99
92) Benzo(g,h,i)perylene	17.06	276	45631	2.321	ng	99

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

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