

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030419\
 Data File : BF112817.D
 Acq On : 4 Mar 2019 12:03
 Operator : JU/SJ
 Sample : K1669-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-031-B

Manual Integrations
 APPROVED

Sohil
 3/5/2019 11:00:31 AM

Quant Time: Mar 05 00:58:12 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	186654	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	664776	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	310492	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	550936	20.00	ng	0.00
76) Chrysene-d12	13.87	240	454993	20.00	ng	0.00
87) Perylene-d12	15.29	264	485843	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	1084049	90.34	ng	0.01
7) Phenol-d6	6.38	99	1408439	93.44	ng	0.01
23) Nitrobenzene-d5	7.30	82	867601	78.09	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	350949	106.47	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1441470	74.44	ng	0.00
79) Terphenyl-d14	12.83	244	1243072	52.96	ng	0.00
Target Compounds						
10) Phenol	6.39	94	64257	3.653	ng	85
50) Dimethylphthalate	9.49	163	103160	4.506	ng	98
71) Phenanthrene	11.27	178	63284	2.309	ng	99
75) Fluoranthene	12.46	202	118421	4.032	ng	95
78) Pyrene	12.68	202	118930	3.101	ng	98
81) Benzo(a)anthracene	13.87	228	70953	2.250	ng	95
83) Chrysene	13.90	228	70306	2.276	ng	96
88) Benzo(b)fluoranthene	14.89	252	102141m	3.250	ng	
90) Benzo(a)pyrene	15.23	252	74699	2.687	ng	94

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

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