

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071619\
 Data File : BF115611.D
 Acq On : 16 Jul 2019 21:31
 Operator : HP/JU
 Sample : K3623-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-071519-A

Manual Integrations
 APPROVED

mohammad
 7/17/2019 12:06:49 PM

Quant Time: Jul 17 05:57:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	203311	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	716966	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	327682	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	618915	20.00	ng	0.00
76) Chrysene-d12	14.04	240	521219	20.00	ng	0.00
87) Perylene-d12	15.52	264	404816	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	440201	35.42	ng	0.00
7) Phenol-d6	6.50	99	549918	34.87	ng	-0.01
23) Nitrobenzene-d5	7.44	82	330164	26.78	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	128554	33.61	ng	-0.01
45) 2-Fluorobiphenyl	9.24	172	606973	32.42	ng	-0.01
79) Terphenyl-d14	12.99	244	668983	23.68	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	76992	3.123	ng	Qvalue 97
71) Phenanthrene	11.43	178	203077	6.401	ng	97
72) Anthracene	11.48	178	70198	2.141	ng	98
75) Fluoranthene	12.62	202	585707	17.471	ng	98
78) Pyrene	12.84	202	582135	13.183	ng	97
81) Benzo(a)anthracene	14.03	228	242642	7.066	ng	99
83) Chrysene	14.07	228	202981	5.889	ng	96
86) Indeno(1,2,3-cd)pyrene	17.00	276	89067	3.041	ng	97
88) Benzo(b)fluoranthene	15.09	252	224222m	8.602	ng	
89) Benzo(k)fluoranthene	15.12	252	62019m	2.669	ng	
90) Benzo(a)pyrene	15.46	252	153802	6.865	ng	# 92
92) Benzo(a,h,i)perylene	17.44	276	86231	4.396	ng	# 92

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

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