

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072219\
 Data File : BF115716.D
 Acq On : 22 Jul 2019 18:20
 Operator : HP/JU
 Sample : K3622-11 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-071919-B

Manual Integrations
 APPROVED

mohammad
 7/23/2019 4:24:45 PM

Quant Time: Jul 23 04:36:44 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	6.88	152	148229	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	571561	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	279237	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	431799	20.00	ng	0.00
76) Chrysene-d12	14.04	240	361727	20.00	ng	0.00
87) Perylene-d12	15.52	264	336070	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	193485	21.35	ng	-0.01
7) Phenol-d6	6.50	99	259106	22.53	ng	-0.01
23) Nitrobenzene-d5	7.43	82	149800	15.24	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	55971	17.17	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	285811	17.91	ng	-0.02
79) Terphenyl-d14	12.99	244	253753	12.94	ng	-0.01
Target Compounds						
71) Phenanthrene	11.43	178	71052	3.210	ng	97
75) Fluoranthene	12.62	202	111133	4.751	ng	99
78) Pyrene	12.84	202	115394	3.765	ng	96
81) Benzo(a)anthracene	14.03	228	50458	2.117	ng	97
88) Benzo(b)fluoranthene	15.09	252	45037m	2.081	ng	

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

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