

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061119\
 Data File : BF114943.D
 Acq On : 11 Jun 2019 13:52
 Operator : HP/JU
 Sample : K3160-05 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-060619-C

Manual Integrations
 APPROVED

mohammad
 6/12/2019 9:03:38 AM

Quant Time: Jun 11 14:34:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 16:44:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	240717	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	924766	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	454839	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	826843	20.00	ng	0.00
76) Chrysene-d12	13.78	240	628472	20.00	ng	-0.01
87) Perylene-d12	15.17	264	619662	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	672107	47.25	ng	0.00
7) Phenol-d6	6.30	99	839340	48.92	ng	-0.01
23) Nitrobenzene-d5	7.22	82	489489	32.52	ng	-0.01
42) 2,4,6-Tribromophenol	10.47	330	220255	44.52	ng	-0.01
45) 2-Fluorobiphenyl	9.02	172	1023484	35.71	ng	0.00
79) Terphenyl-d14	12.75	244	950012	30.06	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.41	163	95324	2.788	ng	98
71) Phenanthrene	11.18	178	270455	6.742	ng	96
75) Fluoranthene	12.36	202	298618	7.197	ng	85
78) Pyrene	12.59	202	331432	6.055	ng	96
81) Benzo(a)anthracene	13.77	228	166186	3.576	ng	96
83) Chrysene	13.80	228	161587	3.463	ng	96
88) Benzo(b)fluoranthene	14.79	252	204103m	5.059	ng	
90) Benzo(a)pyrene	15.12	252	133426	3.808	ng	# 94

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

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