

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
 Data File : BF117863.D
 Acq On : 19 Nov 2019 16:58
 Operator : JU
 Sample : K5865-15
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3-(16-18)

Manual Integrations
 APPROVED

mohammad
 11/20/2019 1:49:21 PM

Quant Time: Nov 20 00:58:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	214255	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	848816	20.00	ng	-0.01
39) Acenaphthene-d10	9.96	164	451897	20.00	ng	-0.01
64) Phenanthrene-d10	11.46	188	750000	20.00	ng	-0.01
76) Chrysene-d12	14.10	240	445453	20.00	ng	-0.01
87) Perylene-d12	15.62	264	466853	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	445441	36.80	ng	-0.01
7) Phenol-d6	6.53	99	1523723	104.37	ng	-0.01
23) Nitrobenzene-d5	7.47	82	1017090	71.23	ng	-0.02
42) 2,4,6-Tribromophenol	10.75	330	8813	1.84	ng	-0.02
45) 2-Fluorobiphenyl	9.28	172	1935498	76.84	ng	-0.01
79) Terphenyl-d14	13.04	244	1791675	73.51	ng	-0.01
Target Compounds						
10) Phenol	6.55	94	73919	4.872	ng	Qvalue 98
50) Dimethylphthalate	9.67	163	220704	6.965	ng	97
71) Phenanthrene	11.48	178	245588	6.513	ng	97
78) Pyrene	12.90	202	183783	5.105	ng	98
81) Benzo(a)anthracene	14.09	228	81018	2.752	ng	96
83) Chrysene	14.13	228	67187	2.355	ng	99
88) Benzo(b)fluoranthene	15.16	252	80418m	2.651	ng	
90) Benzo(a)pyrene	15.54	252	55983	2.099	ng	# 95

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

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