

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112119\
 Data File : BF117905.D
 Acq On : 21 Nov 2019 12:06
 Operator : JU
 Sample : K5927-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-3

Manual Integrations
 APPROVED

mohammad
 11/22/2019 1:20:49 PM

Quant Time: Nov 21 15:35:01 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	154576	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	616051	20.00	ng	-0.01
39) Acenaphthene-d10	9.96	164	302694	20.00	ng	-0.01
64) Phenanthrene-d10	11.46	188	677970	20.00	ng	-0.01
76) Chrysene-d12	14.10	240	541713	20.00	ng	-0.01
87) Perylene-d12	15.62	264	524790	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	887908	101.68	ng	0.00
7) Phenol-d6	6.54	99	1146907	108.89	ng	0.00
23) Nitrobenzene-d5	7.47	82	703361	67.87	ng	-0.02
42) 2,4,6-Tribromophenol	10.76	330	351388	109.65	ng	-0.01
45) 2-Fluorobiphenyl	9.28	172	1424760	84.44	ng	-0.01
79) Terphenyl-d14	13.05	244	1777998	59.99	ng	0.00
Target Compounds						
10) Phenol	6.55	94	28594	2.612	ng	79
50) Dimethylphthalate	9.67	163	248548	11.709	ng	97
71) Phenanthrene	11.48	178	94144	2.762	ng	96
78) Pyrene	12.90	202	127213	2.906	ng	97
81) Benzo(a)anthracene	14.09	228	77560	2.167	ng	99
88) Benzo(b)fluoranthene	15.16	252	79945m	2.345	ng	

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

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