

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
 Data File : BF117621.D
 Acq On : 30 Oct 2019 13:46
 Operator : JU
 Sample : K5539-19
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 165-84-(0-1)

Manual Integrations
 APPROVED

mohammad
 10/31/2019 3:10:34 PM

Quant Time: Oct 31 02:04:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	35258	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	141686	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	65897	20.00	ng	-0.01
64) Phenanthrene-d10	11.25	188	96749	20.00	ng	0.00
76) Chrysene-d12	13.89	240	78569	20.00	ng	0.00
87) Perylene-d12	15.33	264	82202	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	206432	87.10	ng	0.00
7) Phenol-d6	6.38	99	295618	92.87	ng	0.00
23) Nitrobenzene-d5	7.29	82	177506	61.85	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	43164	75.70	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	267496	57.20	ng	-0.01
79) Terphenyl-d14	12.83	244	194003	40.28	ng	0.00
Target Compounds						
10) Phenol	6.40	94	12917	3.976	ng	94
50) Dimethylphthalate	9.48	163	34529	7.350	ng	99
71) Phenanthrene	11.27	178	12217	2.175	ng	96
75) Fluoranthene	12.47	202	18731	3.331	ng	92
78) Pyrene	12.70	202	15998	2.371	ng	99
81) Benzo(a)anthracene	13.89	228	11313	2.135	ng	97
83) Chrysene	13.92	228	11040	2.125	ng	97
88) Benzo(b)fluoranthene	14.93	252	11390m	2.346	ng	

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

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