

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
 Data File : BF117322.D
 Acq On : 3 Oct 2019 18:00
 Operator : JU
 Sample : K5095-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMPOSITE-5

Quant Time: Oct 04 06:44:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	56982	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	229986	20.00	ng	-0.02
39) Acenaphthene-d10	9.84	164	130966	20.00	ng	-0.02
64) Phenanthrene-d10	11.33	188	174462	20.00	ng	0.00
76) Chrysene-d12	13.97	240	90729	20.00	ng	0.00
87) Perylene-d12	15.43	264	85258	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	295999	81.10	ng	0.00
7) Phenol-d6	6.45	99	290655	61.62	ng	0.00
23) Nitrobenzene-d5	7.37	82	309626	70.74	ng	-0.02
42) 2,4,6-Tribromophenol	10.64	330	114377	104.49	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	580918	69.35	ng	-0.01
79) Terphenyl-d14	12.91	244	520911	107.32	ng	0.00

Target Compounds				Qvalue		
10) Phenol	6.46	94	58160	11.185	ng	95
13) 1,4-Dichlorobenzene	6.83	146	13089	2.903	ng	100
14) 1,2-Dichlorobenzene	6.98	146	24866	5.923	ng	97
17) 2-Methylphenol	7.07	107	19619	5.945	ng	# 69
20) 3+4-Methylphenols	7.23	107	46930	11.418	ng	# 48
30) 1,2,4-Trichlorobenzene	8.03	180	18213	5.464	ng	96
31) Naphthalene	8.11	128	144435	13.058	ng	99
32) Benzoic acid	7.87	122	14914	12.380	ng	89
37) 2-Methylnaphthalene	8.80	142	23370	3.138	ng	97
38) 1-Methylnaphthalene	8.90	142	14223m	2.039	ng	
50) Dimethylphthalate	9.55	163	109161	11.772	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.18	232	3991	2.127	ng	# 57
70) Pentachlorophenol	11.15	266	3861	4.227	ng	# 70
74) Di-n-butylphthalate	11.88	149	46815	4.097	ng	97

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

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