

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041019\
 Data File : BF113698.D
 Acq On : 10 Apr 2019 16:00
 Operator : JU/SJ
 Sample : K2316-06RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20190174-AMND-COMP-CS-1RE

Manual Integrations
 APPROVED

Sohil
 4/11/2019 2:39:36 PM

Quant Time: Apr 11 02:16:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	117899	20.00	ng	-0.01
21) Naphthalene-d8	7.96	136	435066	20.00	ng	-0.01
39) Acenaphthene-d10	9.71	164	204138	20.00	ng	-0.01
64) Phenanthrene-d10	11.19	188	373033	20.00	ng	-0.01
76) Chrysene-d12	13.83	240	286175	20.00	ng	-0.02
87) Perylene-d12	15.23	264	304469	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	226222	32.72	ng	0.00
7) Phenol-d6	6.33	99	843364	98.97	ng	0.00
23) Nitrobenzene-d5	7.25	82	586798	79.00	ng	-0.01
42) 2,4,6-Tribromophenol	10.50	330	47109	19.21	ng	-0.02
45) 2-Fluorobiphenyl	9.03	172	826595	65.19	ng	-0.02
79) Terphenyl-d14	12.77	244	773767	55.05	ng	-0.02
Target Compounds						
10) Phenol	6.35	94	22967	2.360	ng	90
50) Dimethylphthalate	9.43	163	158921	10.561	ng	100
71) Phenanthrene	11.22	178	69468	3.748	ng	98
75) Fluoranthene	12.40	202	98476	4.958	ng	98
78) Pyrene	12.63	202	83340	3.827	ng	97
81) Benzo(a)anthracene	13.82	228	36930	2.237	ng	97
83) Chrysene	13.85	228	37798	2.259	ng	97
84) Bis(2-ethylhexyl)phthalate	13.80	149	89453	8.347	ng	97
88) Benzo(b)fluoranthene	14.84	252	49326m	2.647	ng	
90) Benzo(a)pyrene	15.18	252	33954	2.050	ng	# 93

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

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