

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
 Data File : BF111639.D
 Acq On : 20 Dec 2018 15:07
 Operator : JU/SJ
 Sample : J6481-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HOUSE-WALL

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:46:39 PM

Quant Time: Dec 21 06:35:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	204698	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	751987	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	346479	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	536441	20.00	ng	0.00
76) Chrysene-d12	14.04	240	499652	20.00	ng	0.00
87) Perylene-d12	15.52	264	420118	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1093152	91.03	ng	0.00
7) Phenol-d6	6.52	99	1436197	93.97	ng	0.00
23) Nitrobenzene-d5	7.45	82	876432	67.84	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	342590	83.68	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1615792	67.97	ng	0.00
79) Terphenyl-d14	13.00	244	1263401	47.95	ng	0.00
Target Compounds						
10) Phenol	6.53	94	79841	4.630	ng	98
20) 3+4-Methylphenols	7.29	107	43123	3.204	ng	# 56
50) Dimethylphthalate	9.64	163	203698	8.040	ng	97
75) Fluoranthene	12.62	202	127394	4.422	ng	95
78) Pyrene	12.85	202	117679	3.335	ng	94
81) Benzo(a)anthracene	14.03	228	66606	2.336	ng	97
83) Chrysene	14.07	228	72761	2.596	ng	95
84) Bis(2-ethylhexyl)phthalate	14.03	149	51399	2.837	ng	# 98
88) Benzo(b)fluoranthene	15.09	252	89717m	3.654	ng	
90) Benzo(a)pyrene	15.46	252	55161m	2.473	ng	

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

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