

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
 Data File : BF111646.D
 Acq On : 20 Dec 2018 18:42
 Operator : JU/SJ
 Sample : J6346-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DW-10

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:46:48 PM

Quant Time: Dec 21 06:47:21 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	6.89	152	192429	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	721699	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	349916	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	594265	20.00	ng	0.00
76) Chrysene-d12	14.04	240	438737	20.00	ng	0.00
87) Perylene-d12	15.52	264	444786	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	897791	79.53	ng	0.00
7) Phenol-d6	6.52	99	1168864	81.36	ng	0.00
23) Nitrobenzene-d5	7.45	82	717016	57.83	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	294174	71.15	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1404932	58.52	ng	0.00
79) Terphenyl-d14	13.00	244	1099390	47.52	ng	0.00
Target Compounds						
10) Phenol	6.53	94	80376	4.958	ng	92
50) Dimethylphthalate	9.64	163	108181	4.228	ng	97
71) Phenanthrene	11.43	178	94261	2.991	ng	95
75) Fluoranthene	12.62	202	271371	8.503	ng	94
78) Pyrene	12.85	202	200332	6.466	ng	96
81) Benzo(a)anthracene	14.03	228	69227	2.765	ng	97
83) Chrysene	14.07	228	105686	4.295	ng	98
86) Indeno(1,2,3-cd)pyrene	17.00	276	51511	2.462	ng	# 87
88) Benzo(b)fluoranthene	15.09	252	161013m	6.194	ng	
90) Benzo(a)pyrene	15.46	252	79783	3.378	ng	96
92) Benzo(g,h,i)perylene	17.44	276	49004	2.427	ng	# 88

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

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