

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081418\
 Data File : BF108151.D
 Acq On : 14 Aug 2018 18:07
 Operator : JU/SJ
 Sample : J4460-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-080918-SW-002

Manual Integrations
 APPROVED

Sohil
 8/15/2018 3:38:41 PM

Quant Time: Aug 14 19:16:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	172894	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	598767	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	246214	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	430647	20.00	ng	0.00
75) Chrysene-d12	14.22	240	421675	20.00	ng	0.00
86) Perylene-d12	15.78	264	271147	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	818459	85.70	ng	0.02
7) Phenol-d6	6.64	99	1088911	85.81	ng	0.00
23) Nitrobenzene-d5	7.58	82	707845	65.97	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	199843	81.37	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	1082408	70.58	ng	0.00
78) Terphenyl-d14	13.15	244	985218	59.41	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.77	163	119125	6.946	ng	Qvalue # 98
70) Phenanthrene	11.58	178	76859	3.784	ng	98
73) Di-n-butylphthalate	12.10	149	97010	4.630	ng	100
74) Fluoranthene	12.78	202	149031	6.723	ng	96
77) Pyrene	13.01	202	133663	5.007	ng	98
80) Benzo(a)anthracene	14.21	228	90064	3.722	ng	96
82) Chrysene	14.24	228	84369	3.803	ng	97
87) Benzo(b)fluoranthene	15.31	252	90733m	5.600	ng	
89) Benzo(a)pyrene	15.71	252	57156	3.939	ng	95
91) Benzo(g,h,i)perylene	17.89	276	30349	2.567	ng	# 90

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

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