

Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
 Data File : BF097930.D
 Acq On : 22 Aug 2017 2:49
 Operator : SJ/JU
 Sample : I4898-01 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-08-081817-A

Manual Integrations
 APPROVED

Sohil
 8/22/2017 5:45:26 PM

Quant Time: Aug 22 06:57:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	109099	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	382487	20.00	ng	0.00
38) Acenaphthene-d10	9.79	164	142867	20.00	ng	0.00
63) Phenanthrene-d10	11.27	188	257076	20.00	ng	0.00
75) Chrysene-d12	13.90	240	227801	20.00	ng	0.00
86) Perylene-d12	15.32	264	157523	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	406393	56.66	ng	0.00
7) Phenol-d6	6.38	99	510517	52.39	ng	0.00
23) Nitrobenzene-d5	7.31	82	288962	41.04	ng	-0.01
41) 2,4,6-Tribromophenol	10.57	330	64588	52.10	ng	0.00
44) 2-Fluorobiphenyl	9.11	172	360704	37.09	ng	0.00
78) Terphenyl-d14	12.85	244	283600	25.96	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.50	163	85379	8.176	ng	99
70) Phenanthrene	11.29	178	28231	2.061	ng	99
74) Fluoranthene	12.47	202	48898	3.420	ng	98
77) Pyrene	12.70	202	51166	2.746	ng	98
87) Benzo(b)fluoranthene	14.92	252	21124m	2.127	ng	

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

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