

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097042.D
 Acq On : 25 Jul 2017 21:21
 Operator : SJ/JU
 Sample : I4362-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FRAC-TANK

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:54:28 PM

Quant Time: Jul 26 12:20:03 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	160755	20.00	ng	0.00
21) Naphthalene-d8	7.79	136	362133	20.00	ng	0.00
38) Acenaphthene-d10	9.53	164	264235	20.00	ng	0.00
63) Phenanthrene-d10	11.00	188	400690	20.00	ng	0.00
75) Chrysene-d12	13.62	240	261664	20.00	ng	0.00
86) Perylene-d12	14.97	264	213359	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	453794	42.55	ng	0.14
7) Phenol-d6	6.16	99	348504	28.63	ng	0.00
23) Nitrobenzene-d5	7.07	82	640434	103.75	ng	0.00
41) 2,4,6-Tribromophenol	10.32	330	188527	90.30	ng	0.00
44) 2-Fluorobiphenyl	8.86	172	1020332	65.53	ng	0.00
78) Terphenyl-d14	12.59	244	730193	56.90	ng	0.00
Target Compounds						
27) 2,4-Dimethylphenol	7.50	122	64589	11.257	ng	86
31) Naphthalene	7.85	128	10048123	588.624	ng	98
32) Benzoic acid	7.62	122	146678	34.235	ng	67
37) 2-Methylnaphthalene	8.52	142	2441107	225.856	ng	97
51) Acenaphthene	9.56	154	52877	3.521	ng	97
57) Fluorene	10.07	166	88143	5.863	ng	97
70) Phenanthrene	11.03	178	91062	4.242	ng	99

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

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