

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071917\
 Data File : BF096882.D
 Acq On : 19 Jul 2017 20:56
 Operator : SJ/JU
 Sample : I4280-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-071817-A

Manual Integrations
 APPROVED

Sohil
 7/20/2017 3:05:02 PM

Quant Time: Jul 20 05:01:24 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 13:33:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.59	152	113362	20.00	ng	-0.01
21) Naphthalene-d8	7.87	136	423325	20.00	ng	-0.02
38) Acenaphthene-d10	9.62	164	173751	20.00	ng	-0.02
63) Phenanthrene-d10	11.09	188	258884	20.00	ng	-0.01
75) Chrysene-d12	13.72	240	197998	20.00	ng	-0.01
86) Perylene-d12	15.09	264	173597	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	687022	91.12	ng	0.00
7) Phenol-d6	6.24	99	894568	98.87	ng	-0.01
23) Nitrobenzene-d5	7.15	82	455737	66.68	ng	-0.02
41) 2,4,6-Tribromophenol	10.40	330	127865	86.21	ng	-0.02
44) 2-Fluorobiphenyl	8.95	172	791447	71.97	ng	-0.01
78) Terphenyl-d14	12.68	244	499812	43.78	ng	-0.01
Target Compounds						
10) Phenol	6.25	94	36744	3.593	ng	93
49) Dimethylphthalate	9.34	163	127034	9.710	ng	100
74) Fluoranthene	12.30	202	32101	2.383	ng	96
85) Indeno(1,2,3-cd)pyrene	16.37	276	8190	4.030	ng	# 75
87) Benzo(b)fluoranthene	14.72	252	23444m	2.239	ng	

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

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