

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121517\
 Data File : BF101460.D
 Acq On : 15 Dec 2017 23:02
 Operator : SJ/JU
 Sample : I6888-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-122-2(0-5)

Manual Integrations
 APPROVED

Sohil
 12/18/2017 2:04:27 PM

Quant Time: Dec 16 00:33:42 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 15 17:36:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	171885	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	687259	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	282708	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	467297	20.00	ng	0.00
75) Chrysene-d12	14.00	240	334307	20.00	ng	0.00
86) Perylene-d12	15.47	264	301728	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	964146	83.55	ng	0.01
7) Phenol-d6	6.46	99	1164095	81.06	ng	0.00
23) Nitrobenzene-d5	7.39	82	727464	58.92	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	209849	77.03	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1220905	66.99	ng	0.00
78) Terphenyl-d14	12.93	244	895319	64.56	ng	0.00
Target Compounds						
10) Phenol	6.47	94	58929	3.600	ng	89
49) Dimethylphthalate	9.57	163	103612	4.556	ng	99
74) Fluoranthene	12.56	202	66464	2.437	ng	98
77) Pyrene	12.80	202	64297	2.563	ng	97
87) Benzo(b)fluoranthene	15.04	252	50828m	2.764	ng	

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

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