

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\
 Data File : BF102579.D
 Acq On : 29 Jan 2018 1:06
 Operator : SJ/JU
 Sample : J1325-01RE
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-4632RE

Manual Integrations
 APPROVED

Sohil
 1/29/2018 5:48:42 PM

Quant Time: Jan 29 06:17:43 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 27 20:48:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	146165	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	548255	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	201315	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	292846	20.00	ng	0.00
75) Chrysene-d12	13.88	240	279569	20.00	ng	0.00
86) Perylene-d12	15.31	264	212116	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	974423	101.08	ng	0.01
7) Phenol-d6	6.36	99	1148536	98.83	ng	0.00
23) Nitrobenzene-d5	7.29	82	676186	70.88	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	145464	72.92	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1073120	78.38	ng	0.00
78) Terphenyl-d14	12.82	244	781463	63.01	ng	0.00
Target Compounds						
10) Phenol	6.37	94	46566	3.670	ng	85
49) Dimethylphthalate	9.47	163	145668	8.754	ng	100
70) Phenanthrene	11.26	178	120784	7.078	ng	96
74) Fluoranthene	12.45	202	237642	13.656	ng	97
77) Pyrene	12.68	202	215590	9.974	ng	99
80) Benzo(a)anthracene	13.87	228	105008m	6.101	ng	
82) Chrysene	13.90	228	115332m	6.598	ng	
85) Indeno(1,2,3-cd)pyrene	16.71	276	32994	2.286	ng	96
87) Benzo(b)fluoranthene	14.90	252	104737m	7.618	ng	
88) Benzo(k)fluoranthene	14.92	252	41048m	3.160	ng	
89) Benzo(a)pyrene	15.25	252	69435	5.600	ng	# 94
91) Benzo(a,h,i)perylene	17.14	276	33127	3.340	ng	92

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