

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
 Data File : BF111477.D
 Acq On : 15 Dec 2018 19:11
 Operator : JU/SJ
 Sample : J6371-02 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EB07-1-2

Manual Integrations
 APPROVED

Sohil
 12/17/2018 3:12:21 PM

Quant Time: Dec 16 01:48:20 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	178390	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	697445	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	277971	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	400875	20.00	ng	0.00
76) Chrysene-d12	13.96	240	321372	20.00	ng	0.00
87) Perylene-d12	15.40	264	228657	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	255711	23.85	ng	0.00
7) Phenol-d6	6.43	99	738138	51.76	ng	0.00
23) Nitrobenzene-d5	7.35	82	428890	36.62	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	27146	10.90	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	766074	42.57	ng	0.00
79) Terphenyl-d14	12.90	244	546599	39.94	ng	0.00
Target Compounds						
10) Phenol	6.45	94	46486	2.926	ng	91
50) Dimethylphthalate	9.55	163	58968	2.946	ng	# 97
71) Phenanthrene	11.34	178	327534	15.856	ng	93
72) Anthracene	11.39	178	82149	3.889	ng	96
75) Fluoranthene	12.53	202	727805	36.016	ng	97
78) Pyrene	12.76	202	671805	29.650	ng	96
81) Benzo(a)anthracene	13.94	228	338981	19.399	ng	98
83) Chrysene	13.98	228	324165	17.610	ng	97
86) Indeno(1,2,3-cd)pyrene	16.82	276	119420	8.989	ng	# 90
88) Benzo(b)fluoranthene	14.98	252	305424m	22.911	ng	
89) Benzo(k)fluoranthene	15.00	252	99381m	7.802	ng	
90) Benzo(a)pyrene	15.33	252	205216	17.071	ng	98
91) Dibenzo(a,h)anthracene	16.82	278	30223m	3.187	ng	
92) Benzo(a,h,i)perylene	17.24	276	125731	13.507	ng	97

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

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