

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111439.D
 Acq On : 15 Dec 2018 00:40
 Operator : JU/SJ
 Sample : J6374-01 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-029-A

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:37:09 PM

Quant Time: Dec 15 03:04:42 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	247840	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	958483	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	291260	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	439724	20.00	ng	0.00
76) Chrysene-d12	13.95	240	451321	20.00	ng	0.00
87) Perylene-d12	15.39	264	310692	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	640784	43.02	ng	0.00
7) Phenol-d6	6.43	99	1070388	54.03	ng	0.00
23) Nitrobenzene-d5	7.35	82	618384	38.42	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	105938	40.60	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1029974	54.62	ng	0.00
79) Terphenyl-d14	12.90	244	726985	37.83	ng	0.00
Target Compounds						
10) Phenol	6.45	94	64366	2.917	ng	92
50) Dimethylphthalate	9.55	163	62877	2.998	ng	97
71) Phenanthrene	11.33	178	110845	4.892	ng	93
75) Fluoranthene	12.52	202	193794	8.743	ng	98
78) Pyrene	12.75	202	178088	5.597	ng	96
81) Benzo(a)anthracene	13.94	228	85043	3.465	ng	99
83) Chrysene	13.97	228	78766	3.047	ng	97
88) Benzo(b)fluoranthene	14.98	252	71912m	3.970	ng	
90) Benzo(a)pyrene	15.33	252	50240	3.076	ng	# 82
92) Benzo(g,h,i)perylene	17.23	276	27396	2.166	ng	93

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

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