

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071320\
 Data File : BF120851.D
 Acq On : 13 Jul 2020 14:14
 Operator : JU/CG
 Sample : L3184-06
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-070920

Quant Time: Jul 13 14:54:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 14:16:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	196727	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	736121	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	381283	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	705964	20.00	ng	0.00
76) Chrysene-d12	13.99	240	564840	20.00	ng	-0.01
86) Perylene-d12	15.45	264	465212	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1392976	109.30	ng	0.00
7) Phenol-d6	6.46	99	1493445	91.76	ng	0.00
23) Nitrobenzene-d5	7.39	82	1224204	96.35	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	516292	134.78	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1964639	77.96	ng	0.00
79) Terphenyl-d14	12.94	244	2244727	77.81	ng	0.00
Target Compounds						
32) Benzoic acid	7.79	122	1838	4.441	ng	90
50) Dimethylphthalate	9.58	163	299266	10.760	ng	100
87) Indeno(1,2,3-cd)pyrene	16.88	276	85885	2.878	ng	98
91) Dibenzo(a,h)anthracene	16.90	278	68545	2.756	ng	98
92) Benzo(g,h,i)perylene	17.31	276	78479	3.453	ng	99

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

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