

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071521\
 Data File : BF124586.D
 Acq On : 15 Jul 2021 23:07
 Operator : JU/CG
 Sample : M3032-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-07-071421

Manual Integrations
 APPROVED

mohammad
 7/16/2021 10:55:46 AM

Quant Time: Jul 16 02:52:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 07:49:42 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.922	152	4934	20.000	ng	0.00
21) Naphthalene-d8	8.204	136	19441	20.000	ng	0.00
39) Acenaphthene-d10	9.963	164	9754	20.000	ng	0.00
64) Phenanthrene-d10	11.451	188	14502	20.000	ng	0.00
76) Chrysene-d12	14.092	240	10572	20.000	ng	0.00
86) Perylene-d12	15.586	264	10916	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	29142	102.332	ng	0.00
7) Phenol-d6	6.539	99	33427	96.489	ng	0.00
23) Nitrobenzene-d5	7.486	82	20985	68.141	ng	0.00
42) 2,4,6-Tribromophenol	10.751	330	6787	89.415	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	49977	74.716	ng	0.00
79) Terphenyl-d14	13.039	244	40986	65.100	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	11.474	178	3780	4.778	ng	98
75) Fluoranthene	12.662	202	6839	8.568	ng	96
78) Pyrene	12.892	202	6767	7.704	ng	97
81) Benzo(a)anthracene	14.080	228	3362	4.795	ng	99
83) Chrysene	14.115	228	3245m	4.831	ng	
88) Benzo(b)fluoranthene	15.145	252	6326m	8.074	ng	
90) Benzo(a)pyrene	15.521	252	3677	5.685	ng	98

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

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