

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012717.D
 Acq On : 23 Nov 2022 18:57
 Operator : CG/JU
 Sample : N5749-23
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-84

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/25/2022
 Supervised By : Jagrut Upadhyay 11/25/2022

Quant Time: Nov 24 02:32:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	369929	20.000	ng	0.00
21) Naphthalene-d8	10.852	136	1363881	20.000	ng	0.00
39) Acenaphthene-d10	14.663	164	740078	20.000	ng	0.00
64) Phenanthrene-d10	17.422	188	1445213	20.000	ng	0.00
76) Chrysene-d12	21.504	240	1408356	20.000	ng	0.00
86) Perylene-d12	24.027	264	1595731	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2982716	174.678	ng	0.00
7) Phenol-d6	7.181	99	3352528	136.903	ng	0.00
23) Nitrobenzene-d5	9.199	82	2088842	85.415	ng	0.00
42) 2,4,6-Tribromophenol	16.151	330	1254671	131.093	ng	0.00
45) 2-Fluorobiphenyl	13.293	172	4841145	93.193	ng	0.00
79) Terphenyl-d14	19.969	244	7158846	108.035	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.899	128	441816	5.969	ng	98
32) Benzoic acid	10.046	122	25746m	5.246	ng	
74) Di-n-butylphthalate	18.369	149	6634	5.779	ng	98

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

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