

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032222\
 Data File : BP009485.D
 Acq On : 22 Mar 2022 13:33
 Operator : CG/JU
 Sample : N1709-07DL 40X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PPSP-B13-25.5-26.0DL

Quant Time: Mar 23 01:20:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 18:26:53 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 03/23/2022
 Supervised By : Jagrut Upadhyay 03/23/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	458159	20.000	ng	0.00
21) Naphthalene-d8	10.587	136	1621795	20.000	ng	0.00
39) Acenaphthene-d10	14.434	164	921854	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	1774285	20.000	ng	0.00
76) Chrysene-d12	21.304	240	1845763	20.000	ng	-0.01
86) Perylene-d12	23.710	264	2135784	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.399	112	45376	1.729	ng	0.00
7) Phenol-d6	6.981	99	59019	1.663	ng	-0.01
23) Nitrobenzene-d5	8.946	82	53277m	1.746	ng	-0.01
42) 2,4,6-Tribromophenol	15.933	330	11671	0.767	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	85180	1.281	ng	0.00
79) Terphenyl-d14	19.769	244	132666	1.290	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	10.634	128	870777	10.424	ng	96
37) 2-Methylnaphthalene	12.251	142	2342953	39.973	ng	99
38) 1-Methylnaphthalene	12.469	142	1327812	23.254	ng	99
46) 1,1'-Biphenyl	13.263	154	361550	5.257	ng	99
58) Fluorene	15.486	166	228238	3.460	ng	# 86
71) Phenanthrene	17.239	178	759156	7.993	ng	98

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

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