

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
 Data File : BP016440.D
 Acq On : 29 Jul 2023 00:50
 Operator : MA/JU
 Sample : 03770-11
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HOWE-2-LP

Quant Time: Jul 29 03:48:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 01:48:25 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/29/2023
 Supervised By :mohammad ahmed 07/29/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	452454	20.000	ng	0.00
21) Naphthalene-d8	10.922	136	1772008	20.000	ng	0.00
39) Acenaphthene-d10	14.728	164	976769	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	2069334	20.000	ng	0.00
76) Chrysene-d12	21.580	240	1570214	20.000	ng	#-0.01
86) Perylene-d12	24.174	264	1513398	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.605	112	3671709	132.391	ng	0.00
7) Phenol-d6	7.228	99	4260280	112.071	ng	0.00
23) Nitrobenzene-d5	9.287	82	3008574	95.042	ng	0.00
42) 2,4,6-Tribromophenol	16.222	330	1516313	105.522	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	6230587	91.451	ng	0.00
79) Terphenyl-d14	20.033	244	8703224	117.085	ng	-0.01
Target Compounds						
32) Benzoic acid	10.116	122	31924m	7.765	ng	Qvalue

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

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