

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
 Data File : BF137728.D
 Acq On : 24 Apr 2024 20:30
 Operator : CG\JU
 Sample : P2226-01RE
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GROUND-WATERRE

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/26/2024
 Supervised By :mohammad ahmed 04/27/2024

Quant Time: Apr 25 04:20:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 03:57:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.063	152	285671	20.000	ng	0.00
21) Naphthalene-d8	8.345	136	1136938	20.000	ng	0.00
39) Acenaphthene-d10	10.110	164	604286	20.000	ng	0.00
64) Phenanthrene-d10	11.604	188	901316	20.000	ng	0.00
76) Chrysene-d12	14.251	240	485789	20.000	ng	0.00
86) Perylene-d12	15.827	264	604041	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	825845	43.914	ng	0.00
7) Phenol-d6	6.669	99	690471	28.912	ng	0.00
23) Nitrobenzene-d5	7.622	82	1770823	82.679	ng	0.00
42) 2,4,6-Tribromophenol	10.898	330	505687	82.931	ng	0.00
45) 2-Fluorobiphenyl	9.422	172	3193504	78.965	ng	0.00
79) Terphenyl-d14	13.192	244	2532756	83.918	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.681	94	578528	24.249	ng	97
32) Benzoic acid	8.110	122	694826m	57.280	ng	

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

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