

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092024\
 Data File : BF139508.D
 Acq On : 20 Sep 2024 18:20
 Operator : RC/JU
 Sample : P4103-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-14-5-7.5A

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/21/2024
 Supervised By :mohammad ahmed 09/22/2024

Quant Time: Sep 20 23:21:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 14 02:58:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	74397	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	256441	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	119498	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	212737	20.000	ng	0.00
76) Chrysene-d12	14.033	240	125388	20.000	ng	# 0.00
86) Perylene-d12	15.510	264	91312	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	547931	120.148	ng	0.02
7) Phenol-d6	6.510	99	662312	110.645	ng	0.00
23) Nitrobenzene-d5	7.440	82	498516	91.389	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	151836	119.349	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	779918	91.537	ng	0.00
79) Terphenyl-d14	12.986	244	813837	106.535	ng	0.00
Target Compounds						
70) Pentachlorophenol	11.204	266	9131m	6.150	ng	Qvalue

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

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