Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139003.D
Acq On : 14 Aug 2024 19:46
Operator : RC/JU
Sample : P3476-05RE
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MUL06RE

Quant Time: Aug 15 01:12:32 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 30 17:50:01 2024
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	6.834	152	30464	20.000	ng		-0.01
21) Naphthalene-d8	8.116	136	138299	20.000	ng		-0.01
39) Acenaphthene-d10	9.869	164	51885	20.000	ng		-0.01
64) Phenanthrene-d10	11.357	188	95379	20.000	ng		-0.01
76) Chrysene-d12	13.998	240	63161	20.000	ng		0.00
86) Perylene-d12	15.463	264	47717	20.000	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.475	112	118723	60.159	ng		0.00
7) Phenol-d6	6.492	99	90975	34.335	ng		0.00
23) Nitrobenzene-d5	7.404	82	221124	78.172	ng		0.00
42) 2,4,6-Tribromophenol	10.663	330	62022	145.931	ng		0.00
45) 2-Fluorobiphenyl	9.186	172	305681	88.520	ng		-0.02
79) Terphenyl-d14	12.939	244	406711	107.811	ng		0.00
Target Compounds							Qvalue
10) Phenol	6.504	94	13106	4.698	ng		91
15) Benzyl Alcohol	6.992	79	67760	35.482	ng		95
20) 3+4-Methylphenols	7.263	107	29583	13.448	ng	#	48
22) Acetophenone	7.245	105	17003	5.021	ng	#	94
32) Benzoic acid	7.922	122	35499	30.479	ng		98

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

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