

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092024\
 Data File : BN034067.D
 Acq On : 20 Sep 2024 02:57
 Operator : JU/RC
 Sample : P3957-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW305D-20240910

Quant Time: Sep 20 03:50:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.438	152	5360	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	15498	0.400	ng	0.00
13) Acenaphthene-d10	14.083	164	9061	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	21444	0.400	ng	# 0.00
29) Chrysene-d12	21.053	240	16206	0.400	ng	0.00
35) Perylene-d12	23.177	264	14678	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	2134	0.140	ng	0.00
5) Phenol-d6	6.636	99	1294	0.073	ng	0.00
8) Nitrobenzene-d5	8.578	82	3634	0.307	ng	0.00
11) 2-Methylnaphthalene-d10	11.791	152	8153	0.356	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	1429	0.348	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	13665	0.371	ng	0.00
27) Fluoranthene-d10	18.880	212	24802	0.458	ng	0.00
31) Terphenyl-d14	19.498	244	16700	0.516	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.075	88	4214	0.629	ng	# 74
24) Pentachlorophenol	16.498	266	5873	1.314	ng	100

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

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