

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082124\
 Data File : BP021602.D
 Acq On : 21 Aug 2024 18:08
 Operator : RC/JU
 Sample : P3617-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 Field Blank

Quant Time: Aug 22 00:20:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 14 00:00:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	252809	20.000	ng	-0.01
21) Naphthalene-d8	10.593	136	1002689	20.000	ng	# 0.00
39) Acenaphthene-d10	14.451	164	670558	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	1647270	20.000	ng	0.01
76) Chrysene-d12	21.727	240	1735021	20.000	ng	0.03
86) Perylene-d12	25.180	264	1891478	20.000	ng	0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	1156998	67.956	ng	0.00
7) Phenol-d6	6.963	99	978428	39.395	ng	0.00
23) Nitrobenzene-d5	8.946	82	2122867	109.278	ng	0.00
42) 2,4,6-Tribromophenol	15.975	330	1568727	210.199	ng	0.00
45) 2-Fluorobiphenyl	13.063	172	3972812	90.419	ng	0.00
79) Terphenyl-d14	19.992	244	10278440	115.089	ng	0.02

Target Compounds	Qvalue

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

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