

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042925\
 Data File : BP024462.D
 Acq On : 29 Apr 2025 16:41
 Operator : RC/JU
 Sample : Q1871-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-A

Quant Time: Apr 29 17:03:42 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 18 12:04:48 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	157409	20.000	ng	-0.01
21) Naphthalene-d8	10.481	136	619557	20.000	ng	#-0.02
39) Acenaphthene-d10	14.351	164	392799	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	853964	20.000	ng	# 0.00
76) Chrysene-d12	21.598	240	985729	20.000	ng	0.00
86) Perylene-d12	24.939	264	903330	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	1141102	119.864	ng	-0.01
7) Phenol-d6	6.899	99	1260131	96.669	ng	-0.01
23) Nitrobenzene-d5	8.857	82	949622	87.399	ng	-0.01
42) 2,4,6-Tribromophenol	15.869	330	917152	168.762	ng	0.00
45) 2-Fluorobiphenyl	12.957	172	2058562	80.242	ng	0.00
79) Terphenyl-d14	19.880	244	4601001	94.082	ng	0.00
Target Compounds						
32) Benzoic acid	9.787	122	110832m	14.401	ng	Qvalue

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

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