

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021924\
 Data File : BP019767.D
 Acq On : 19 Feb 2024 17:28
 Operator : MA/JU
 Sample : P1461-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-1

Quant Time: Feb 19 17:53:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	58866	20.000	ng	-0.02
21) Naphthalene-d8	10.699	136	208653	20.000	ng	-0.02
39) Acenaphthene-d10	14.534	164	145032	20.000	ng	-0.02
64) Phenanthrene-d10	17.292	188	370279	20.000	ng	-0.01
76) Chrysene-d12	21.386	240	412121	20.000	ng	# 0.00
86) Perylene-d12	23.804	264	449954	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	451941	123.755	ng	-0.01
7) Phenol-d6	7.069	99	538785	109.418	ng	-0.02
23) Nitrobenzene-d5	9.069	82	433545	90.035	ng	-0.02
42) 2,4,6-Tribromophenol	16.034	330	413963	195.490	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	1020657	86.999	ng	-0.02
79) Terphenyl-d14	19.857	244	2408032	96.091	ng	-0.01
Target Compounds						Qvalue
83) Chrysene	21.422	228	65974	2.395	ng	98
84) Bis(2-ethylhexyl)phtha...	21.304	149	35532	2.148	ng	98
85) Di-n-octyl phthalate	22.245	149	60122	2.066	ng	# 84
88) Benzo(b)fluoranthene	23.063	252	77058m	2.717	ng	
89) Benzo(k)fluoranthene	23.116	252	74218	2.740	ng	97

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

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