

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021924\
 Data File : BP019768.D
 Acq On : 19 Feb 2024 18:02
 Operator : MA/JU
 Sample : P1505-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-1S

Quant Time: Feb 19 23:36:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 19 23:35:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	47588	20.000	ng	0.00
21) Naphthalene-d8	10.699	136	167547	20.000	ng	0.00
39) Acenaphthene-d10	14.534	164	108381	20.000	ng	0.00
64) Phenanthrene-d10	17.292	188	274481	20.000	ng	0.00
76) Chrysene-d12	21.380	240	332580	20.000	ng	-0.01
86) Perylene-d12	23.804	264	382285	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.470	112	345050	116.877	ng	0.00
7) Phenol-d6	7.069	99	385321	96.797	ng	0.00
23) Nitrobenzene-d5	9.069	82	336127	86.930	ng	0.00
42) 2,4,6-Tribromophenol	16.028	330	286019	180.746	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	740068	84.415	ng	0.00
79) Terphenyl-d14	19.857	244	1702700	84.195	ng	0.00
Target Compounds						
83) Chrysene	21.422	228	50829	2.287	ng	Qvalue 96
84) Bis(2-ethylhexyl)phtha...	21.304	149	28467	2.133	ng	# 98
88) Benzo(b)fluoranthene	23.069	252	56603	2.349	ng	95
89) Benzo(k)fluoranthene	23.116	252	72820m	3.164	ng	

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

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