

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092521\
 Data File : BN016479.D
 Acq On : 25 Sep 2021 17:32
 Operator : CG/JU
 Sample : M3892-21
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EB-01-092221

Quant Time: Sep 26 04:15:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 03:35:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.670	152	20592	0.400	ng	0.00
7) Naphthalene-d8	10.432	136	88733	0.400	ng	#-0.01
13) Acenaphthene-d10	14.294	164	38254	0.400	ng	0.00
19) Phenanthrene-d10	17.042	188	83678	0.400	ng	0.00
29) Chrysene-d12	21.259	240	67881	0.400	ng	# 0.01
35) Perylene-d12	23.518	264	23615	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	6409	0.139	ng	0.02
5) Phenol-d6	6.849	99	3982	0.073	ng	0.00
8) Nitrobenzene-d5	8.810	82	21464	0.342	ng	0.00
11) 2-Methylnaphthalene-d10	12.029	152	43961	0.334	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	3170	0.347	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	63263	0.394	ng	-0.01
27) Fluoranthene-d10	19.087	212	89487	0.390	ng	0.00
31) Terphenyl-d14	19.698	244	96602	0.612	ng	0.00
Target Compounds						
						Qvalue
33) Chrysene	21.291	228	12999	0.061	ng	98
34) Bis(2-ethylhexyl)phtha...	21.185	149	7640	0.142	ng	96
36) Indeno(1,2,3-cd)pyrene	25.815	276	6301	0.082	ng	95
37) Benzo(b)fluoranthene	22.837	252	10810	0.137	ng	94
38) Benzo(k)fluoranthene	22.882	252	9837	0.129	ng	# 81
40) Dibenzo(a,h)anthracene	25.836	278	6650	0.102	ng	# 88
41) Benzo(g,h,i)perylene	26.513	276	4559	0.068	ng	98

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

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